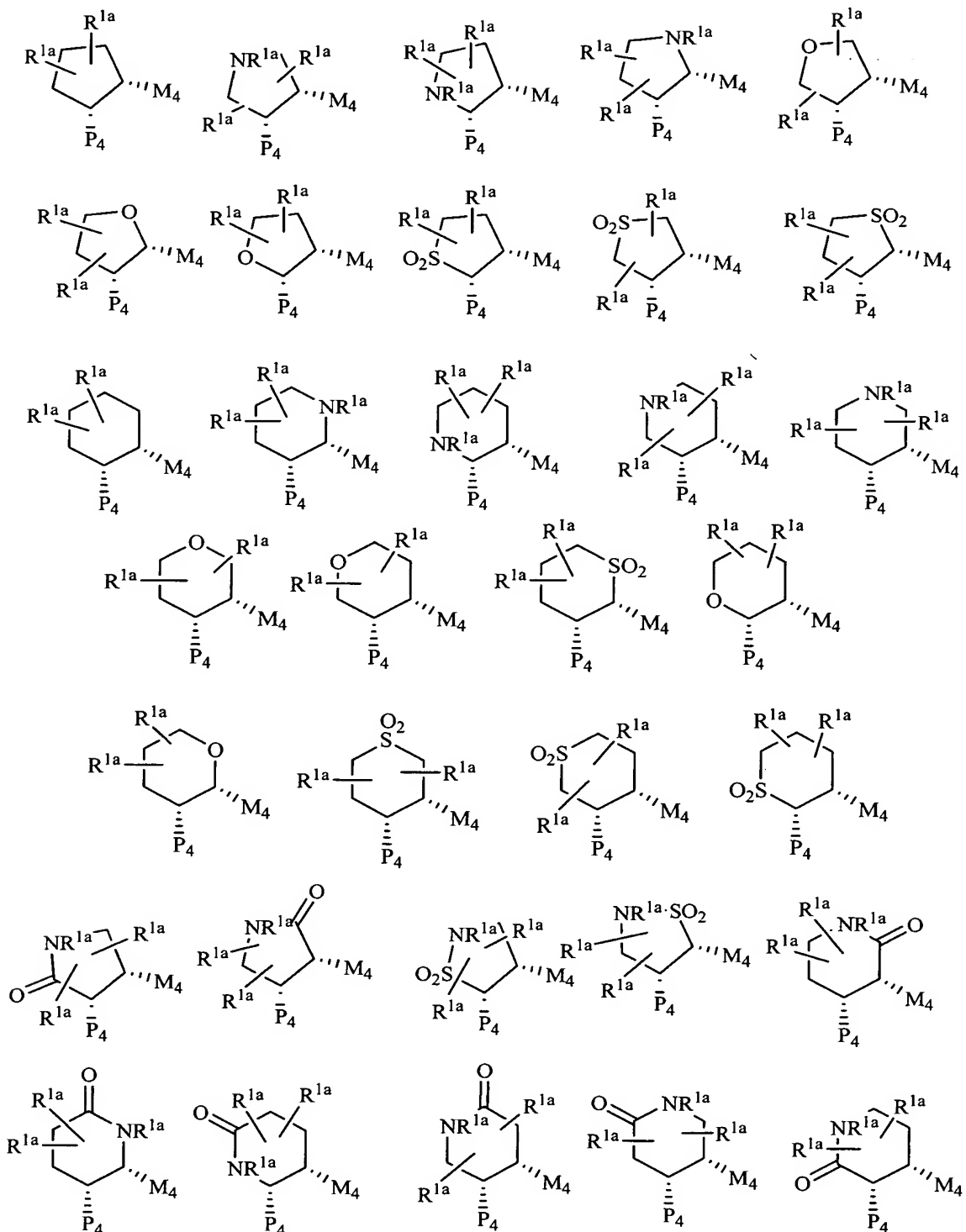
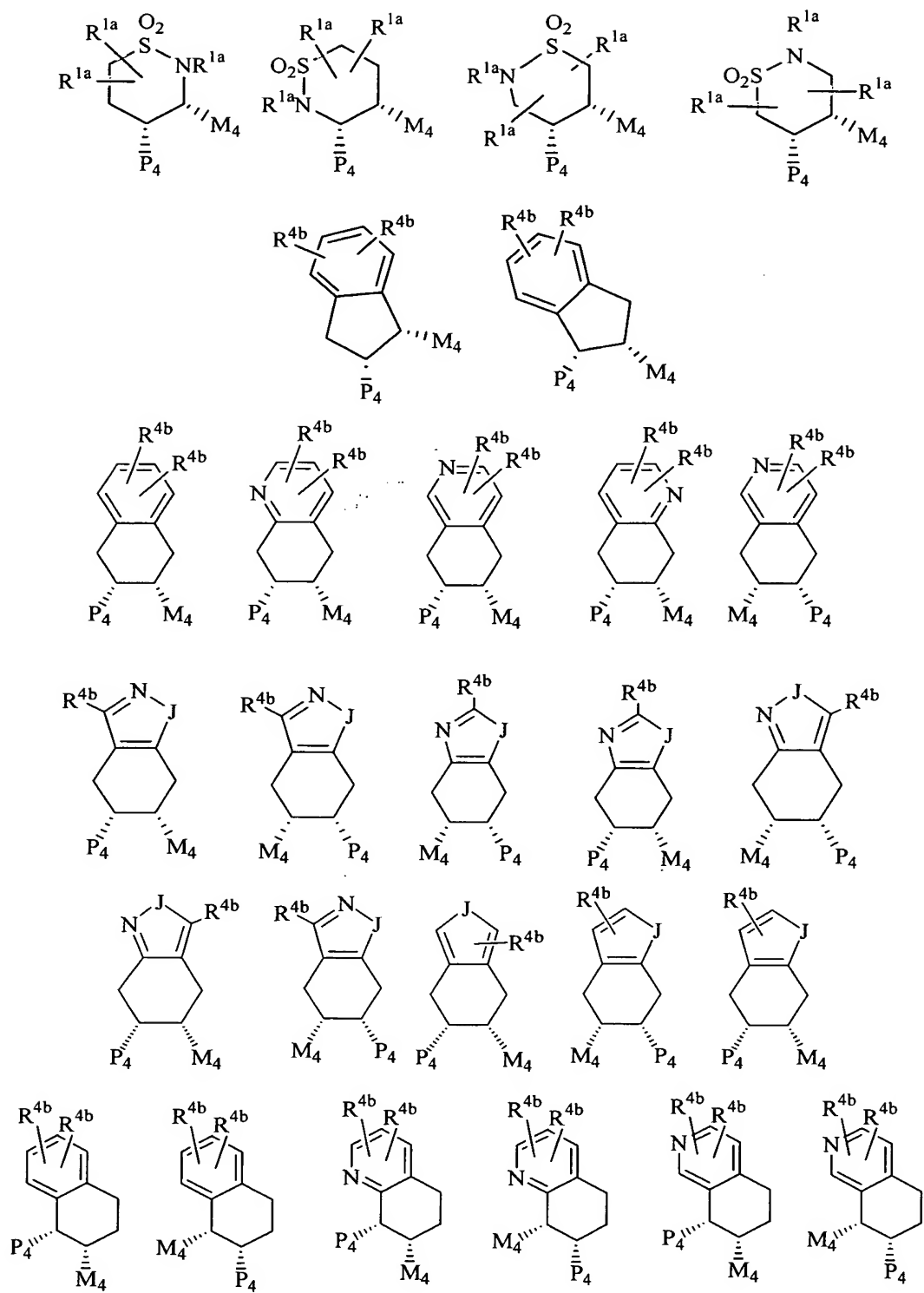
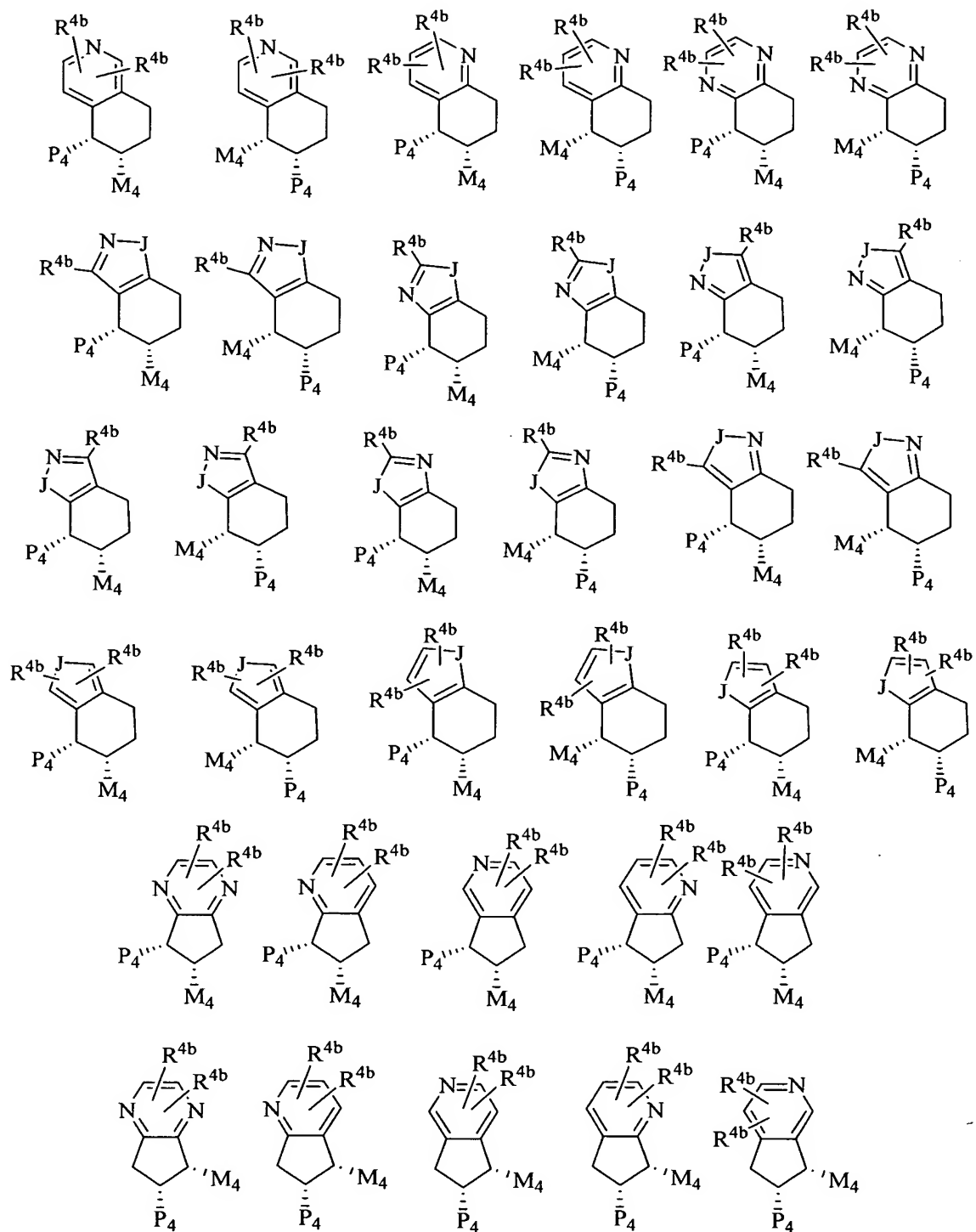


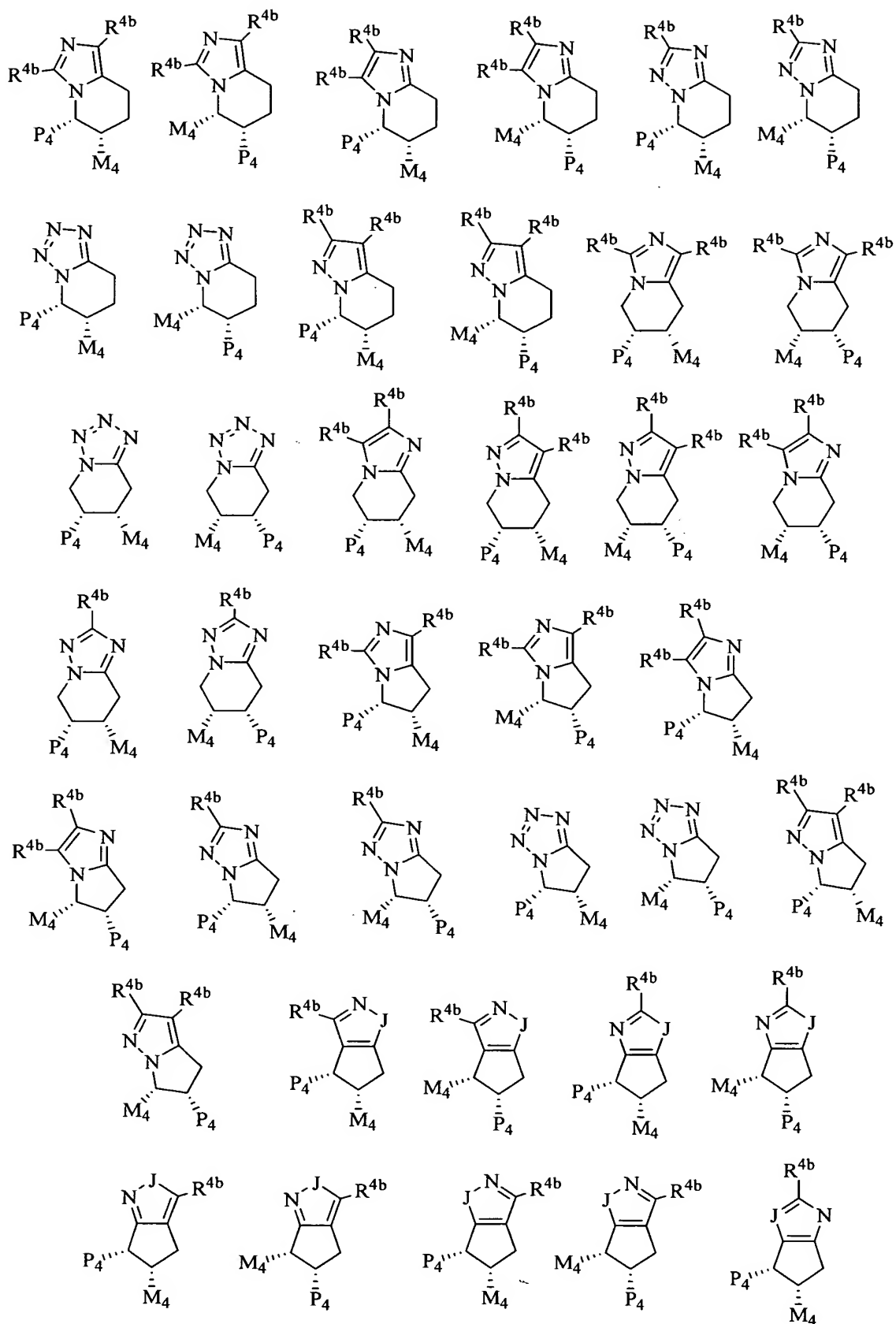
WHAT IS CLAIMED IS:

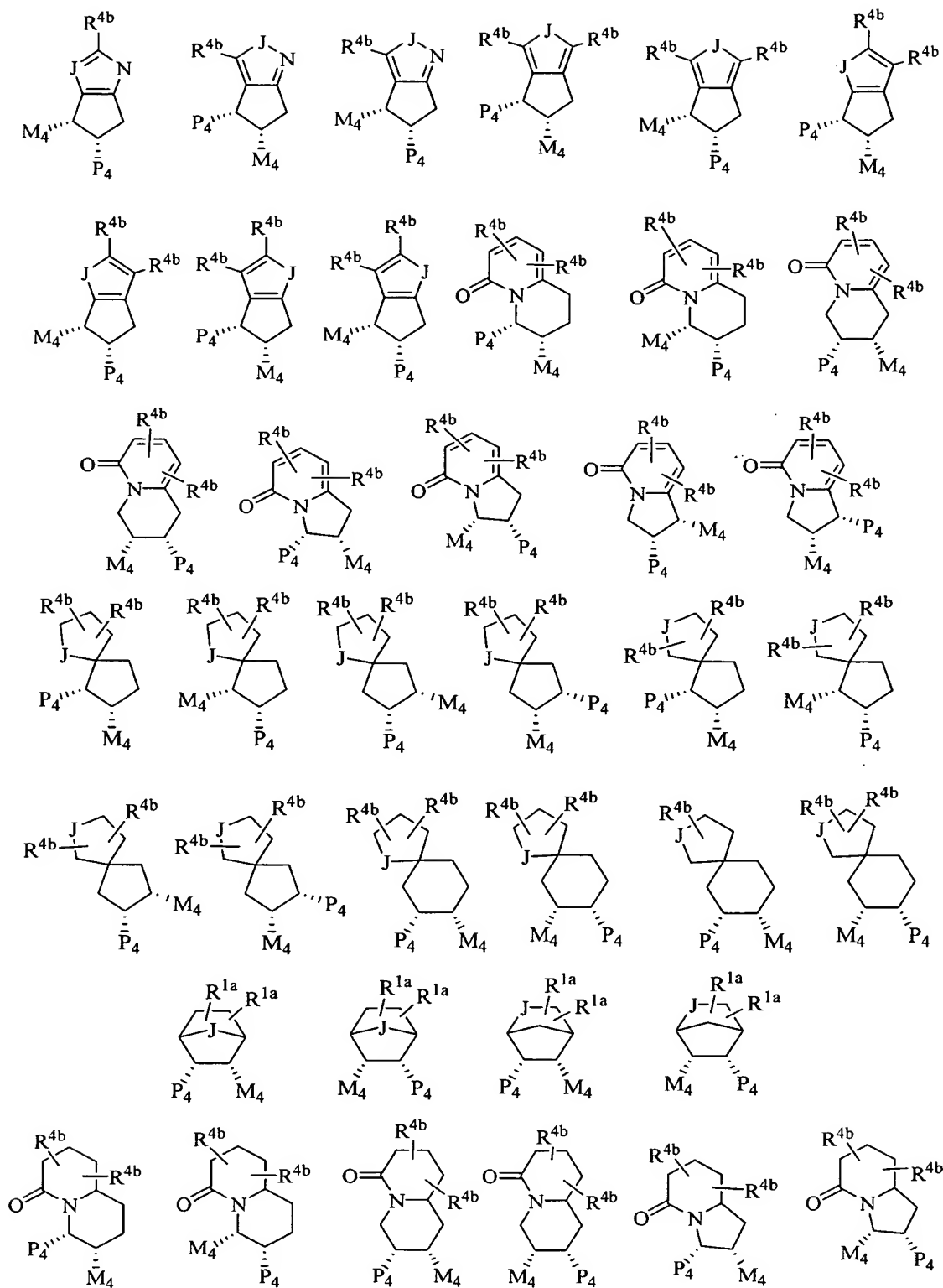
1. A compound selected from:

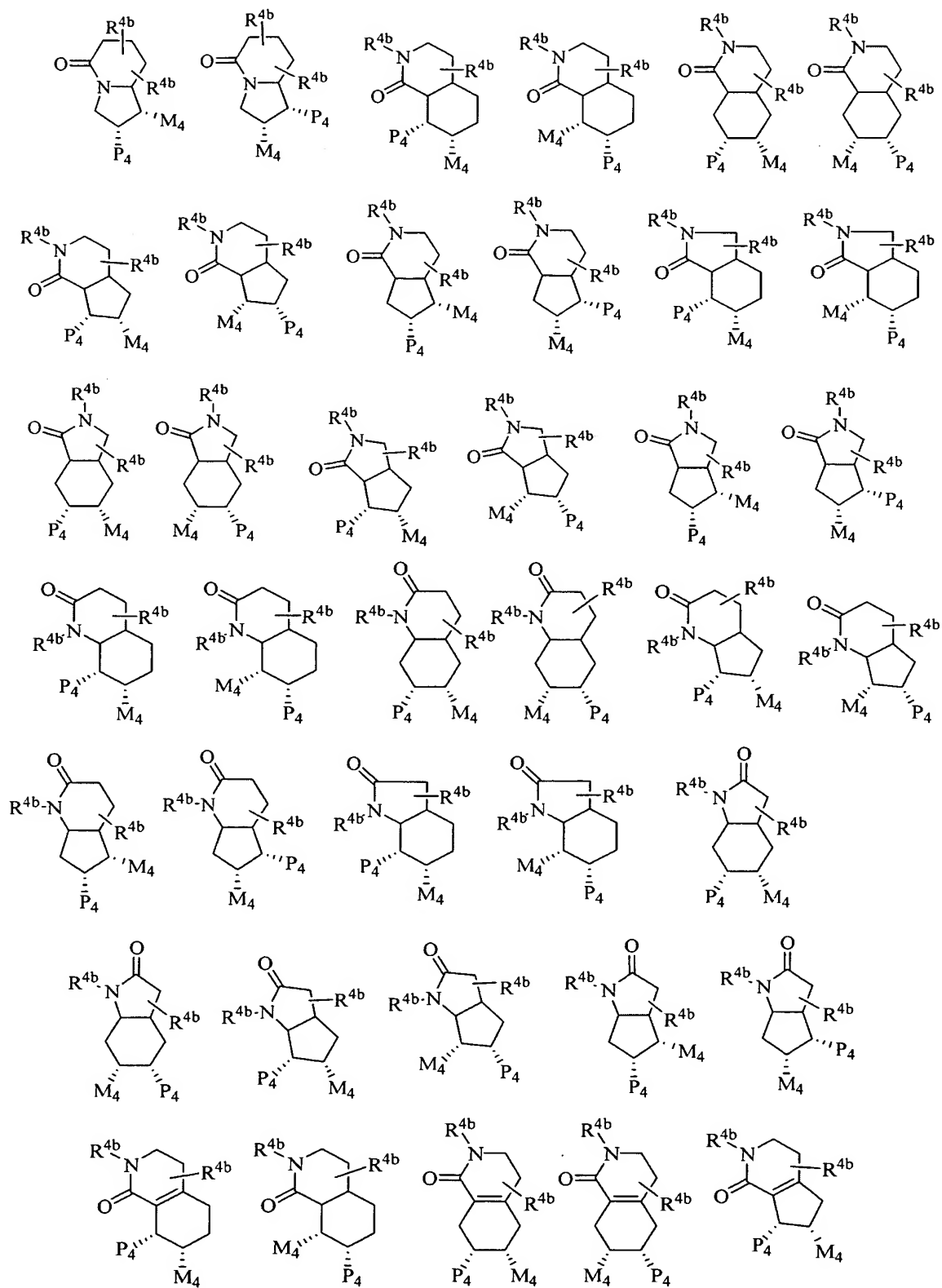


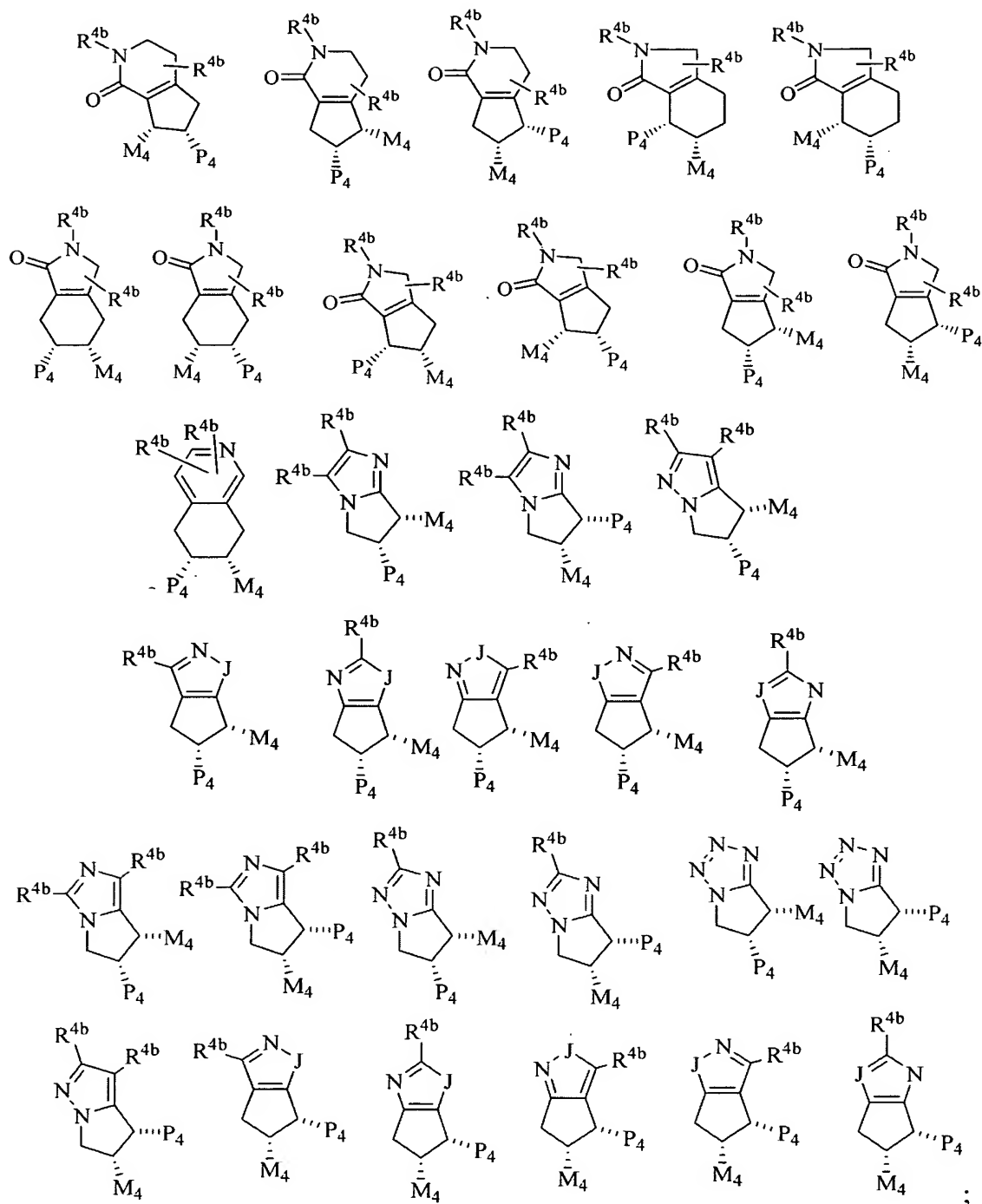






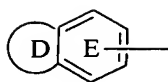




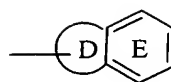


or a stereoisomer or pharmaceutically acceptable salt thereof, wherein:

- 5 J is selected from O, S, S(O)₂, CR^{1a}, and NR^{1a};
 one of P₄ and M₄ is -Z-A-B and the other -G₁-G;
 G is a group of formula IIa or IIb:



IIa



IIb

ring D, including the two atoms of Ring E to which it is attached, is a 5-6
 5 membered ring consisting of carbon atoms and 0-2 heteroatoms selected from the
 group consisting of N, O, and S(O)_p;

ring D is substituted with 0-2 R and there are 0-3 ring double bonds;

E is selected from phenyl, pyridyl, pyrimidyl, pyrazinyl, and pyridazinyl, and
 is substituted with 1-3 R;

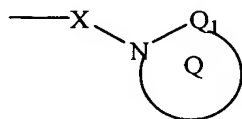
alternatively, ring D is absent and ring E is selected from phenyl, pyridyl,
 10 pyrimidyl, pyrazinyl, pyridazinyl, pyrrolyl, pyrazolyl, imidazolyl, isoxazolyl,
 oxazolyl, triazolyl, thienyl, and thiazolyl, and ring E is substituted with 1-3 R;

alternatively, ring D is absent and ring E is selected from phenyl, pyridyl,
 pyrimidyl, pyrazinyl, pyridazinyl, pyrrolyl, pyrazolyl, imidazolyl, isoxazolyl,
 oxazolyl, triazolyl, thienyl, and thiazolyl, and ring E is substituted with 1 R and with a
 15 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected
 from the group consisting of N, O, and S(O)_p, wherein the 5-6 membered heterocycle
 is substituted with 0-2 carbonyls and 1-3 R and there are 0-3 ring double bonds;

R is selected from H, C₁₋₄ alkyl, F, Cl, Br, I, OH, OCH₃, OCH₂CH₃,
 OCH(CH₃)₂, OCH₂CH₂CH₃, -CN, NH₂, NH(C₁₋₃ alkyl), N(C₁₋₃ alkyl)₂,
 20 C(=NH)NH₂, CH₂NH₂, CH₂NH(C₁₋₃ alkyl), CH₂N(C₁₋₃ alkyl)₂, CH₂CH₂NH₂,
 CH₂CH₂NH(C₁₋₃ alkyl), CH₂CH₂N(C₁₋₃ alkyl)₂, C(=NR⁸)NR⁷R⁹,
 NHC(=NR⁸)NR⁷R⁹, ONHC(=NR⁸)NR⁷R⁹, NR⁸CH(=NR⁷), (CR⁸R⁹)_tC(O)H,
 (CR⁸R⁹)_tC(O)R^{2c}, (CR⁸R⁹)_tNR⁷R⁸, (CR⁸R⁹)_tC(O)NR⁷R⁸, (CR⁸R⁹)_tNR⁷C(O)R⁷,
 (CR⁸R⁹)_tOR³, (CR⁸R⁹)_tS(O)_pNR⁷R⁸, (CR⁸R⁹)_tNR⁷S(O)_pR⁷, (CR⁸R⁹)_tSR³,
 25 (CR⁸R⁹)_tS(O)R³, (CR⁸R⁹)_tS(O)₂R³, and OCF₃, provided that S(O)_pR⁷ and S(O)₂R³
 form other than S(O)₂H or S(O)H;

alternatively, when 2 R groups are attached to adjacent atoms, they combine to
 form methylenedioxy or ethylenedioxy;

A is selected from: C₃₋₁₀ carbocycle substituted with 0-2 R⁴, and 5-12 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p and substituted with 0-2 R⁴;



5 B is ; provided that Z and B are attached to different atoms on A and that the A-X-N moiety forms other than a N-N-N group;

Q₁ is selected from C=O and SO₂;

ring Q is a 4-7 membered monocyclic or tricyclic ring consisting of, in addition to the N-Q₁ group shown, carbon atoms and 0-2 heteroatoms selected from
 10 NR^{4c}, O, and S(O)_p, wherein: 0-2 double bonds are present within the ring and the ring is substituted with 0-2 R^{4a};

alternatively, ring Q is a 4-7 membered ring to which another ring is fused, wherein: the 4-7 membered ring consists of, in addition to the N-Q₁ group shown, carbon atoms and 0-2 heteroatoms selected from NR^{4c}, O, and S(O)_p and 0-1 double
 15 bonds are present within the ring; the fusion ring is phenyl or a 5-6 membered heteroaromatic consisting of carbon atoms and 1-2 heteroatoms selected from NR^{4c}, O, and S(O)_p;

ring Q, which includes the 4-7 membered ring and the fusion ring, is substituted with 0-3 R^{4a};

20 X is absent or is selected from (CR²R^{2a})₁₋₄, C(O), C(O)CR²R^{2a}, CR²R^{2a}C(O), S(O)₂, S(O)₂CR²R^{2a}, CR²R^{2a}S(O)₂, NR²S(O)₂, NR²CR²R^{2a}, and OCR²R^{2a}, wherein the left side of X is attached to ring A;

G₁ is selected from (CR³R^{3a})₁₋₅, (CR³R^{3a})₀₋₂CR³=CR³(CR³R^{3a})₀₋₂, (CR³R^{3a})₀₋₂C≡C(CR³R^{3a})₀₋₂, (CR³R^{3a})_uC(O)(CR³R^{3a})_w,
 25 (CR³R^{3a})_uC(O)O(CR³R^{3a})_w, (CR³R^{3a})_uOC(O)(CR³R^{3a})_w, (CR³R^{3a})_uO(CR³R^{3a})_w, (CR³R^{3a})_uNR^{3b}(CR³R^{3a})_w, (CR³R^{3a})_uNR^{3e}(CR³R^{3a})_w, (CR³R^{3a})_uC(O)NR^{3b}(CR³R^{3a})_w, (CR³R^{3a})_uNR^{3b}C(O)(CR³R^{3a})_w, (CR³R^{3a})_uOC(O)NR^{3b}(CR³R^{3a})_w, (CR³R^{3a})_uNR^{3b}C(O)O(CR³R^{3a})_w,

- $(\text{CR}^3\text{R}^{3a})_u\text{NR}^{3b}\text{C}(\text{O})(\text{CR}^3\text{R}^{3a})_u\text{C}(\text{O})(\text{CR}^3\text{R}^{3a})_w$,
 $(\text{CR}^3\text{R}^{3a})_u\text{C}(\text{O})(\text{CR}^3\text{R}^{3a})_u\text{C}(\text{O})\text{NR}^{3b}(\text{CR}^3\text{R}^{3a})_w$,
 $(\text{CR}^3\text{R}^{3a})_u\text{NR}^{3b}\text{C}(\text{O})(\text{CR}^3\text{R}^{3a})_u\text{C}(\text{O})\text{NR}^{3b}(\text{CR}^3\text{R}^{3a})_w$,
 $(\text{CR}^3\text{R}^{3a})_u\text{S}(\text{O})_2\text{NR}^{3b}\text{C}(\text{O})(\text{CR}^3\text{R}^{3a})_w$, $(\text{CR}^3\text{R}^{3a})_u\text{C}(\text{O})\text{NR}^{3b}\text{S}(\text{O})_2(\text{CR}^3\text{R}^{3a})_w$,
5 $(\text{CR}^3\text{R}^{3a})_u\text{S}(\text{O})_2\text{NR}^{3b}\text{C}(\text{O})\text{NR}^{3b}(\text{CR}^3\text{R}^{3a})_w$,
 $(\text{CR}^3\text{R}^{3a})_u\text{NR}^{3b}\text{C}(\text{O})(\text{CR}^3\text{R}^{3a})_u\text{C}(\text{S})\text{NR}^{3b}(\text{CR}^3\text{R}^{3a})_w$, and
 $(\text{CR}^3\text{R}^{3a})_u\text{NR}^{3b}\text{C}(\text{S})(\text{CR}^3\text{R}^{3a})_u\text{C}(\text{O})\text{NR}^{3b}(\text{CR}^3\text{R}^{3a})_w$, wherein $u+w$ or $u+u+w$ total 0,
1, 2, 3, or 4, and the right side of Z is attached to ring A, provided that Z does not
form an N-S, NCH₂N, NCH₂O, or NCH₂S bond with either group to which it is
10 attached;
 R^{1a} , at each occurrence, is selected from H, $-(\text{CR}^3\text{R}^{3a})_r\text{R}^{1b}$,
 $-(\text{CR}^3\text{R}^{3a})_r\text{CR}^3\text{R}^{1b}\text{R}^{1b}$, $-(\text{CR}^3\text{R}^{3a})_r\text{O}-(\text{CR}^3\text{R}^{3a})_r\text{R}^{1b}$, $-\text{C}_{2-6}$ alkenylene- R^{1b} ,
 $-\text{C}_{2-6}$ alkynylene- R^{1b} , $-(\text{CR}^3\text{R}^{3a})_r\text{C}(=\text{NR}^{1b})\text{NR}^3\text{R}^{1b}$, $\text{NR}^3\text{CR}^3\text{R}^{3a}\text{R}^{1c}$, $\text{OCR}^3\text{R}^{3a}\text{R}^{1c}$,
 $\text{SCR}^3\text{R}^{3a}\text{R}^{1c}$, $\text{NR}^3(\text{CR}^3\text{R}^{3a})_2(\text{CR}^3\text{R}^{3a})_t\text{R}^{1b}$, $\text{C}(\text{O})\text{NR}^2(\text{CR}^3\text{R}^{3a})_2(\text{CR}^3\text{R}^{3a})_t\text{R}^{1b}$,
15 $\text{CO}_2(\text{CR}^3\text{R}^{3a})_2(\text{CR}^3\text{R}^{3a})_t\text{R}^{1b}$, $\text{O}(\text{CR}^3\text{R}^{3a})_2(\text{CR}^3\text{R}^{3a})_t\text{R}^{1b}$, $\text{S}(\text{CR}^3\text{R}^{3a})_2(\text{CR}^3\text{R}^{3a})_t\text{R}^{1b}$,
 $\text{S}(\text{O})_p(\text{CR}^3\text{R}^{3a})_t\text{R}^{1d}$, $\text{O}(\text{CR}^3\text{R}^{3a})_t\text{R}^{1d}$, $\text{NR}^3(\text{CR}^3\text{R}^{3a})_t\text{R}^{1d}$, $\text{OC}(\text{O})\text{NR}^3(\text{CR}^3\text{R}^{3a})_t\text{R}^{1d}$,
 $\text{NR}^3\text{C}(\text{O})\text{NR}^3(\text{CR}^3\text{R}^{3a})_t\text{R}^{1d}$, $\text{NR}^3\text{C}(\text{O})\text{O}(\text{CR}^3\text{R}^{3a})_t\text{R}^{1d}$, and $\text{NR}^3\text{C}(\text{O})(\text{CR}^3\text{R}^{3a})_t\text{R}^{1d}$,
provided that R^{1a} forms other than an N-halo, N-S, O-O, or N-CN bond;
alternatively, when two R^{1a} groups are attached to adjacent atoms, together
20 with the atoms to which they are attached they form a 5-7 membered ring consisting
of: carbon atoms and 0-2 heteroatoms selected from the group consisting of N, O,
and $\text{S}(\text{O})_p$, this ring being substituted with 0-2 R^{4b} and having 0-3 ring double bonds;
 R^{1b} is selected from H, C_{1-3} alkyl, F, Cl, Br, I, -CN, -NO₂, -CHO, $(\text{CF}_2)_r\text{CF}_3$,
 $(\text{CR}^3\text{R}^{3a})_r\text{OR}^2$, NR^2R^{2a} , $\text{C}(\text{O})\text{R}^{2b}$, CO_2R^{2b} , $\text{OC}(\text{O})\text{R}^2$, $(\text{CF}_2)_r\text{CO}_2\text{R}^{2a}$, $\text{S}(\text{O})_p\text{R}^{2b}$,
25 $\text{NR}^2(\text{CH}_2)_t\text{OR}^2$, $\text{C}(=\text{NR}^{2c})\text{NR}^2\text{R}^{2a}$, $\text{NR}^2\text{C}(\text{O})\text{R}^{2b}$, $\text{NR}^2\text{C}(\text{O})\text{NHR}^2$, $\text{NR}^2\text{C}(\text{O})_2\text{R}^{2a}$,
 $\text{OC}(\text{O})\text{NR}^2\text{R}^{2a}$, $\text{C}(\text{O})\text{NR}^2\text{R}^{2a}$, $\text{C}(\text{O})\text{NR}^2\text{R}^{2b}$, $\text{C}(\text{S})\text{NR}^2\text{R}^{2a}$, $\text{C}(\text{O})\text{NR}^2(\text{CH}_2)_t\text{OR}^2$,
 $\text{SO}_2\text{NR}^2\text{R}^{2a}$, $\text{NR}^2\text{SO}_2\text{R}^2$, $\text{C}(\text{O})\text{NR}^2\text{SO}_2\text{R}^2$, C_{3-6} carbocycle substituted with 0-2 R^{4b} ,
and 5-10 membered heterocycle consisting of carbon atoms and from 1-4 heteroatoms

selected from the group consisting of N, O, and S(O)_p, and substituted with 0-2 R^{4b}, provided that R^{1b} forms other than an O-O, N-halo, N-S, or N-CN bond;

R^{1c} is selected from H, CH(CH₂OR²)₂, C(O)R^{2c}, C(O)NR²R^{2a}, S(O)R², S(O)₂R², and SO₂NR²R^{2a};

5 R^{1d} is selected from C₃₋₆ carbocycle substituted with 0-2 R^{4b} and 5-10 membered heterocycle consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-2 R^{4b}, provided that R^{1d} forms other than an N-S bond;

R², at each occurrence, is selected from H, CF₃, C₁₋₆ alkyl, -(CH₂)_r-C₃₋₁₀ carbocycle substituted with 0-2 R^{4b}, and -(CH₂)_r-5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-2 R^{4b};

R^{2a}, at each occurrence, is selected from H, CF₃, C₁₋₆ alkyl, -(CH₂)_r-C₃₋₁₀ carbocycle substituted with 0-2 R^{4b}, and -(CH₂)_r-5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-2 R^{4b};

alternatively, NR²R^{2a} forms a 4, 5, or 6 membered saturated, partially saturated, or unsaturated ring substituted with 0-2 R^{4b} and consisting of: carbon atoms, the nitrogen atom to which R² and R^{2a} are attached, and 0-1 additional heteroatoms selected from the group consisting of N, O, and S(O)_p;

R^{2b}, at each occurrence, is selected from CF₃, C₁₋₄ alkoxy substituted with 0-2 R^{4b}, C₁₋₆ alkyl substituted with 0-3 R^{4b}, -(CH₂)_r-C₃₋₁₃ carbocycle substituted with 0-2 R^{4b}, and -(CH₂)_r-5-10 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-2 R^{4b};

R^{2c}, at each occurrence, is selected from CF₃, OH, C₁₋₄ alkoxy, C₁₋₆ alkyl, -(CH₂)_r-C₃₋₁₀ carbocycle substituted with 0-2 R^{4b}, and -(CH₂)_r-5-10 membered heterocycle containing from 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-2 R^{4b};

R^3 , at each occurrence, is selected from H, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, benzyl, and phenyl;

R^{3a} , at each occurrence, is selected from H, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$,
 5 $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, benzyl, and phenyl;

alternatively, NR^3R^{3a} forms a 5 or 6 membered saturated, partially unsaturated, or unsaturated ring consisting of: carbon atoms, the nitrogen atom to which R^3 and R^{3a} are attached, and 0-1 additional heteroatoms selected from the
 10 group consisting of N, O, and $S(O)_p$;

R^{3b} , at each occurrence, is selected from H, C_{1-6} alkyl substituted with 0-2 R^{1a} , C_{2-6} alkenyl substituted with 0-2 R^{1a} , C_{2-6} alkynyl substituted with 0-2 R^{1a} , $-(C_{0-4}$ alkyl)-5-10 membered carbocycle substituted with 0-3 R^{1a} , and $-(C_{0-4}$ alkyl)-5-10 membered heterocycle substituted with 0-3 R^{1a} and consisting of: carbon
 15 atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{3c} , at each occurrence, is selected from CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, benzyl, and phenyl;

R^{3d} , at each occurrence, is selected from H, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$,
 20 $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, C_{1-4} alkyl-phenyl, and $C(=O)R^{3c}$;

R^{3e} , at each occurrence, is selected from H, SO_2NHR^3 , $SO_2NR^3R^3$, $C(O)R^3$, $C(O)NHR^3$, $C(O)OR^{3f}$, $S(O)R^{3f}$, $S(O)_2R^{3f}$, C_{1-6} alkyl substituted with 0-2 R^{1a} , C_{2-6} alkenyl substituted with 0-2 R^{1a} , C_{2-6} alkynyl substituted with 0-2 R^{1a} , $-(C_{0-4}$ alkyl)-5-10 membered carbocycle substituted with 0-3 R^{1a} , and $-(C_{0-4}$ alkyl)-5-10 membered heterocycle substituted with 0-3 R^{1a} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{3f} , at each occurrence, is selected from: C_{1-6} alkyl substituted with 0-2 R^{1a} , C_{2-6} alkenyl substituted with 0-2 R^{1a} , C_{2-6} alkynyl substituted with 0-2 R^{1a} , $-(C_{0-4}$

alkyl)-5-10 membered carbocycle substituted with 0-3 R^{1a} , and $-(C_{0-4} \text{ alkyl})$ -5-10 membered heterocycle substituted with 0-3 R^{1a} and consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p;

- R^4 , at each occurrence, is selected from H, =O, $(CR^3R^{3a})_rOR^2$, F, Cl, Br, I,
- 5 C_{1-4} alkyl, $(CR^3R^{3a})_rCN$, $(CR^3R^{3a})_rNO_2$, $(CR^3R^{3a})_rNR^2R^{2a}$, $(CR^3R^{3a})_rC(O)R^{2c}$, $(CR^3R^{3a})_rNR^2C(O)R^{2b}$, $(CR^3R^{3a})_rC(O)NR^2R^{2a}$, $(CR^3R^{3a})_rNR^2C(O)NR^2R^{2a}$, $(CR^3R^{3a})_rC(=NR^2)NR^2R^{2a}$, $(CR^3R^{3a})_rC(=NS(O)_2R^5)NR^2R^{2a}$, $(CR^3R^{3a})_rNHC(=NR^2)NR^2R^{2a}$, $(CR^3R^{3a})_rC(O)NHC(=NR^2)NR^2R^{2a}$, $(CR^3R^{3a})_rSO_2NR^2R^{2a}$, $(CR^3R^{3a})_rNR^2SO_2NR^2R^{2a}$, $(CR^3R^{3a})_rNR^2SO_2-C_{1-4}$ alkyl,
- 10 $(CR^3R^{3a})_rNR^2SO_2R^5$, $(CR^3R^{3a})_rS(O)_pR^{5a}$, $(CR^3R^{3a})_r(CF_2)_rCF_3$, $NHCH_2R^{1c}$, OCH_2R^{1c} , SCH_2R^{1c} , $NH(CH_2)_2(CH_2)_tR^{1b}$, $O(CH_2)_2(CH_2)_tR^{1b}$, $S(CH_2)_2(CH_2)_tR^{1b}$, $(CR^3R^{3a})_{r-5-6}$ membered carbocycle substituted with 0-1 R^5 , and a $(CR^3R^{3a})_{r-5-6}$ membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p and substituted with 0-1 R^5 ;
- 15 R^{4a} , at each occurrence, is selected from H, =O, $(CR^3R^{3a})_rOR^2$, $(CR^3R^{3a})_rF$, $(CR^3R^{3a})_rBr$, $(CR^3R^{3a})_rCl$, C_{1-4} alkyl, $(CR^3R^{3a})_rCN$, $(CR^3R^{3a})_rNO_2$, $(CR^3R^{3a})_rNR^2R^{2a}$, $(CR^3R^{3a})_rC(O)R^{2c}$, $(CR^3R^{3a})_rNR^2C(O)R^{2b}$, $(CR^3R^{3a})_rC(O)NR^2R^{2a}$, $(CR^3R^{3a})_rN=CHOR^3$, $(CR^3R^{3a})_rC(O)NH(CH_2)_2NR^2R^{2a}$, $(CR^3R^{3a})_rNR^2C(O)NR^2R^{2a}$, $(CR^3R^{3a})_rC(=NR^2)NR^2R^{2a}$,
- 20 $(CR^3R^{3a})_rNHC(=NR^2)NR^2R^{2a}$, $(CR^3R^{3a})_rSO_2NR^2R^{2a}$, $(CR^3R^{3a})_rNR^2SO_2NR^2R^{2a}$, $(CR^3R^{3a})_rNR^2SO_2-C_{1-4}$ alkyl, $(CR^3R^{3a})_rC(O)NHSO_2-C_{1-4}$ alkyl, $(CR^3R^{3a})_rNR^2SO_2R^5$, $(CR^3R^{3a})_rS(O)_pR^{5a}$, $(CR^3R^{3a})_r(CF_2)_rCF_3$, $(CR^3R^{3a})_{r-5-6}$ membered carbocycle substituted with 0-1 R^5 , and a $(CR^3R^{3a})_{r-5-6}$ membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group
- 25 consisting of N, O, and S(O)_p, and substituted with 0-1 R^5 ;

R^{4b} , at each occurrence, is selected from H, =O, $(CH_2)_rOR^3$, $(CH_2)_rF$, $(CH_2)_rCl$, $(CH_2)_rBr$, $(CH_2)_rI$, C_{1-4} alkyl, $(CH_2)_rCN$, $(CH_2)_rNO_2$, $(CH_2)_rNR^3R^{3a}$, $(CH_2)_rC(O)R^3$, $(CH_2)_rC(O)OR^{3c}$, $(CH_2)_rNR^3C(O)R^{3a}$, $(CH_2)_rC(O)NR^3R^{3a}$, $(CH_2)_rNR^3C(O)NR^3R^{3a}$, $(CH_2)_rC(=NR^3)NR^3R^{3a}$, $(CH_2)_rNR^3C(=NR^3)NR^3R^{3a}$,

$(\text{CH}_2)_r\text{SO}_2\text{NR}^3\text{R}^{3a}$, $(\text{CH}_2)_r\text{NR}^3\text{SO}_2\text{NR}^3\text{R}^{3a}$, $(\text{CH}_2)_r\text{NR}^3\text{SO}_2\text{-C}_{1-4}$ alkyl,
 $(\text{CH}_2)_r\text{NR}^3\text{SO}_2\text{CF}_3$, $(\text{CH}_2)_r\text{NR}^3\text{SO}_2\text{-phenyl}$, $(\text{CH}_2)_r\text{S(O)}_p\text{CF}_3$, $(\text{CH}_2)_r\text{S(O)}_p\text{-C}_{1-4}$
 alkyl, $(\text{CH}_2)_r\text{S(O)}_p\text{-phenyl}$, and $(\text{CH}_2)_r(\text{CF}_2)_r\text{CF}_3$;

- R^{4c} , at each occurrence, is selected from H, C_{1-4} alkyl, $(\text{CR}^3\text{R}^{3a})_{r1}\text{OR}^2$,
- 5 $(\text{CR}^3\text{R}^{3a})_{r1}\text{F}$, $(\text{CR}^3\text{R}^{3a})_{r1}\text{Br}$, $(\text{CR}^3\text{R}^{3a})_{r1}\text{Cl}$, $(\text{CR}^3\text{R}^{3a})_{r1}\text{CN}$, $(\text{CR}^3\text{R}^{3a})_{r1}\text{NO}_2$,
 $(\text{CR}^3\text{R}^{3a})_{r1}\text{NR}^2\text{R}^{2a}$, $(\text{CR}^3\text{R}^{3a})_{r1}\text{C(O)}\text{R}^{2c}$, $(\text{CR}^3\text{R}^{3a})_{r1}\text{NR}^2\text{C(O)}\text{R}^{2b}$,
 $(\text{CR}^3\text{R}^{3a})_{r1}\text{C(O)}\text{NR}^2\text{R}^{2a}$, $(\text{CR}^3\text{R}^{3a})_{r1}\text{N=CHOR}^3$, $(\text{CR}^3\text{R}^{3a})_{r1}\text{C(O)}\text{NH}(\text{CH}_2)_2\text{NR}^2\text{R}^{2a}$,
 $(\text{CR}^3\text{R}^{3a})_{r1}\text{NR}^2\text{C(O)}\text{NR}^2\text{R}^{2a}$, $(\text{CR}^3\text{R}^{3a})_{r1}\text{C(=NR}^2)\text{NR}^2\text{R}^{2a}$,
 $(\text{CR}^3\text{R}^{3a})_{r1}\text{NHC(=NR}^2)\text{NR}^2\text{R}^{2a}$, $(\text{CR}^3\text{R}^{3a})_{r1}\text{SO}_2\text{NR}^2\text{R}^{2a}$,
- 10 $(\text{CR}^3\text{R}^{3a})_{r1}\text{NR}^2\text{SO}_2\text{NR}^2\text{R}^{2a}$, $(\text{CR}^3\text{R}^{3a})_{r1}\text{NR}^2\text{SO}_2\text{-C}_{1-4}$ alkyl,
 $(\text{CR}^3\text{R}^{3a})_{r1}\text{C(O)}\text{NHSO}_2\text{-C}_{1-4}$ alkyl, $(\text{CR}^3\text{R}^{3a})_{r1}\text{NR}^2\text{SO}_2\text{R}^5$, $(\text{CR}^3\text{R}^{3a})_{r1}\text{S(O)}_p\text{R}^{5a}$,
 $(\text{CR}^3\text{R}^{3a})_{r1}(\text{CF}_2)_r\text{CF}_3$, $(\text{CR}^3\text{R}^{3a})_{r1-5-6}$ membered carbocycle substituted with 0-1 R^5 ,
 and a $(\text{CR}^3\text{R}^{3a})_{r1-5-6}$ membered heterocycle consisting of: carbon atoms and 1-4
 heteroatoms selected from the group consisting of N, O, and S(O)_p , and substituted
- 15 with 0-1 R^5 ;

- R^5 , at each occurrence, is selected from H, C_{1-6} alkyl, $=\text{O}$, $(\text{CH}_2)_r\text{OR}^3$, F, Cl,
 Br, I, $-\text{CN}$, NO_2 , $(\text{CH}_2)_r\text{NR}^3\text{R}^{3a}$, $(\text{CH}_2)_r\text{C(O)}\text{R}^3$, $(\text{CH}_2)_r\text{C(O)}\text{OR}^{3c}$,
 $(\text{CH}_2)_r\text{NR}^3\text{C(O)}\text{R}^{3a}$, $(\text{CH}_2)_r\text{C(O)}\text{NR}^3\text{R}^{3a}$, $(\text{CH}_2)_r\text{NR}^3\text{C(O)}\text{NR}^3\text{R}^{3a}$,
 $(\text{CH}_2)_r\text{CH(=NOR}^{3d})$, $(\text{CH}_2)_r\text{C(=NR}^3)\text{NR}^3\text{R}^{3a}$, $(\text{CH}_2)_r\text{NR}^3\text{C(=NR}^3)\text{NR}^3\text{R}^{3a}$,
- 20 $(\text{CH}_2)_r\text{SO}_2\text{NR}^3\text{R}^{3a}$, $(\text{CH}_2)_r\text{NR}^3\text{SO}_2\text{NR}^3\text{R}^{3a}$, $(\text{CH}_2)_r\text{NR}^3\text{SO}_2\text{-C}_{1-4}$ alkyl,
 $(\text{CH}_2)_r\text{NR}^3\text{SO}_2\text{CF}_3$, $(\text{CH}_2)_r\text{NR}^3\text{SO}_2\text{-phenyl}$, $(\text{CH}_2)_r\text{S(O)}_p\text{CF}_3$, $(\text{CH}_2)_r\text{S(O)}_p\text{-C}_{1-4}$
 alkyl, $(\text{CH}_2)_r\text{S(O)}_p\text{-phenyl}$, $(\text{CF}_2)_r\text{CF}_3$, phenyl substituted with 0-2 R^6 , naphthyl
 substituted with 0-2 R^6 , and benzyl substituted with 0-2 R^6 ;

- R^{5a} , at each occurrence, is selected from C_{1-6} alkyl, $(\text{CH}_2)_r\text{OR}^3$,
- 25 $(\text{CH}_2)_r\text{NR}^3\text{R}^{3a}$, $(\text{CH}_2)_r\text{C(O)}\text{R}^3$, $(\text{CH}_2)_r\text{C(O)}\text{OR}^{3c}$, $(\text{CH}_2)_r\text{NR}^3\text{C(O)}\text{R}^{3a}$,
 $(\text{CH}_2)_r\text{C(O)}\text{NR}^3\text{R}^{3a}$, $(\text{CF}_2)_r\text{CF}_3$, phenyl substituted with 0-2 R^6 , naphthyl substituted
 with 0-2 R^6 , and benzyl substituted with 0-2 R^6 , provided that R^{5a} does not form a
 S-N or $\text{S(O)}_p\text{-C(O)}$ bond;

R^6 , at each occurrence, is selected from H, OH, $(CH_2)_rOR^2$, Cl, F, Br, I, C_{1-4} alkyl, -CN, NO_2 , $(CH_2)_rNR^2R^{2a}$, $(CH_2)_rC(O)R^{2b}$, $NR^2C(O)R^{2b}$, $NR^2C(O)NR^2R^{2a}$, $C(=NH)NH_2$, $NHC(=NH)NH_2$, $SO_2NR^2R^{2a}$, $NR^2SO_2NR^2R^{2a}$, and $NR^2SO_2C_{1-4}$ alkyl;

5 R^7 , at each occurrence, is selected from H, OH, C_{1-6} alkyl, C_{1-6} alkyl-C(O)-, C_{1-6} alkyl-O-, $(CH_2)_n$ -phenyl, C_{1-4} alkyl-OC(O)-, C_{6-10} aryl-O-, C_{6-10} aryl-OC(O)-, C_{6-10} aryl-CH₂C(O)-, C_{1-4} alkyl-C(O)O- C_{1-4} alkyl-OC(O)-, C_{6-10} aryl-C(O)O- C_{1-4} alkyl-OC(O)-, C_{1-6} alkyl-NH₂-C(O)-, phenyl-NH₂-C(O)-, and phenyl- C_{1-4} alkyl-C(O)-;

10 R^8 , at each occurrence, is selected from H, C_{1-6} alkyl, and $(CH_2)_n$ -phenyl;
alternatively, NR^7R^8 forms a 5-10 membered heterocyclic ring consisting of carbon atoms and 0-2 additional heteroatoms selected from the group consisting of N, O, and S(O)_p;

R^9 , at each occurrence, is selected from H, C_{1-6} alkyl, and $(CH_2)_n$ -phenyl;

15 n, at each occurrence, is selected from 0, 1, 2, and 3;

p, at each occurrence, is selected from 0, 1, and 2;

r, at each occurrence, is selected from 0, 1, 2, 3, 4, 5, and 6;

r1, at each occurrence, is selected from 1, 2, 3, 4, 5, and 6; and

t, at each occurrence, is selected from 0, 1, 2, and 3.

20

2. A compound according to Claim 1, wherein:

G is selected from the group: 2-aminomethyl-4-chloro-phenyl;

2-aminosulfonyl-4-chloro-phenyl; 2-amido-4-chloro-phenyl;

25 4-chloro-2-methylsulfonyl-phenyl; 2-aminosulfonyl-4-fluoro-phenyl;

2-amido-4-fluoro-phenyl; 4-fluoro-2-methylsulfonyl-phenyl;

2-aminomethyl-4-bromo-phenyl; 2-aminosulfonyl-4-bromo-phenyl;

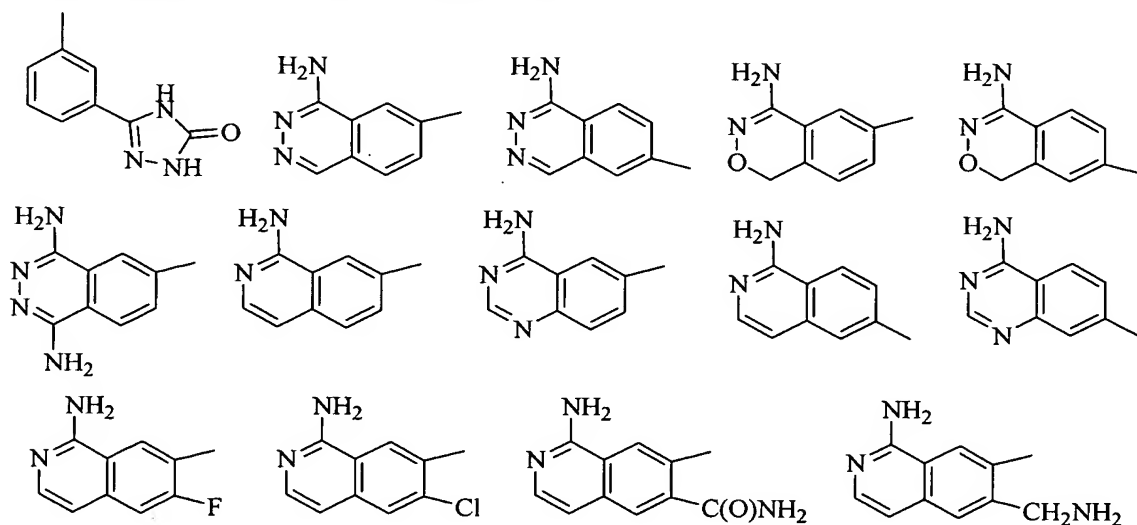
2-amido-4-bromo-phenyl; 4-bromo-2-methylsulfonyl-phenyl;

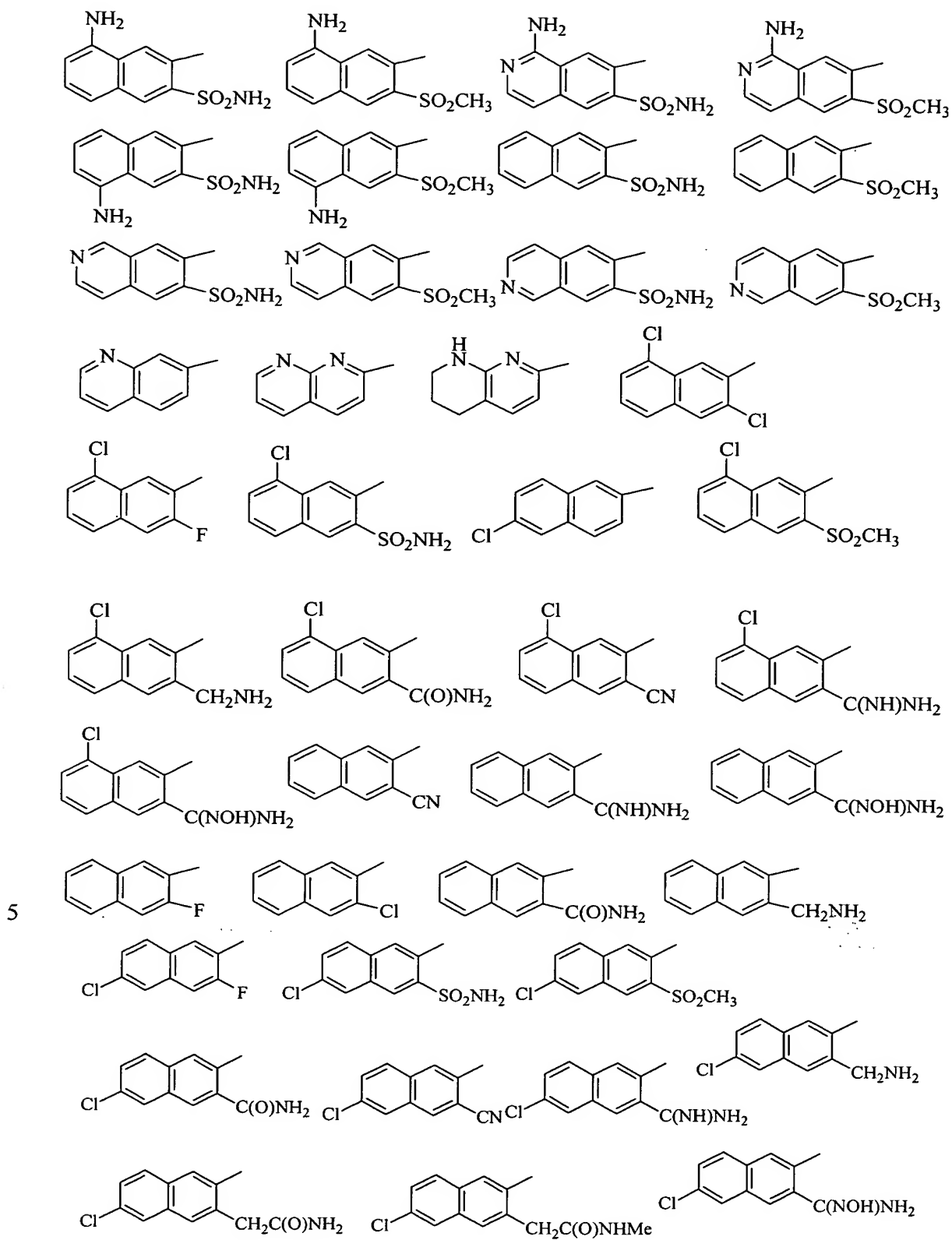
2-aminomethyl-4-methyl-phenyl; 2-aminosulfonyl-4-methyl-phenyl;

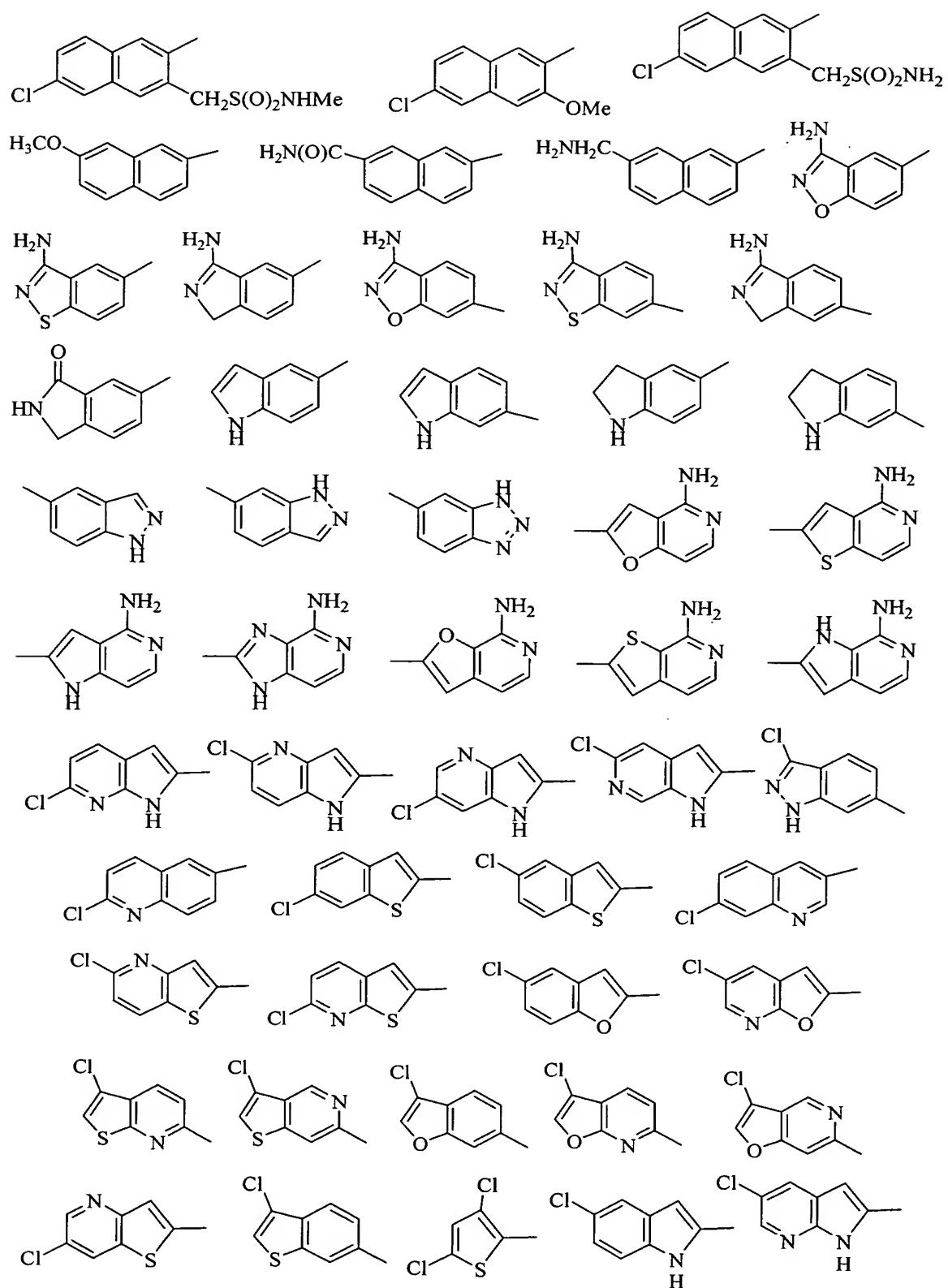
30 2-amido-4-methyl-phenyl; 2-methylsulfonyl-4-methyl-phenyl; 4-fluoro-pyrid-2-yl;

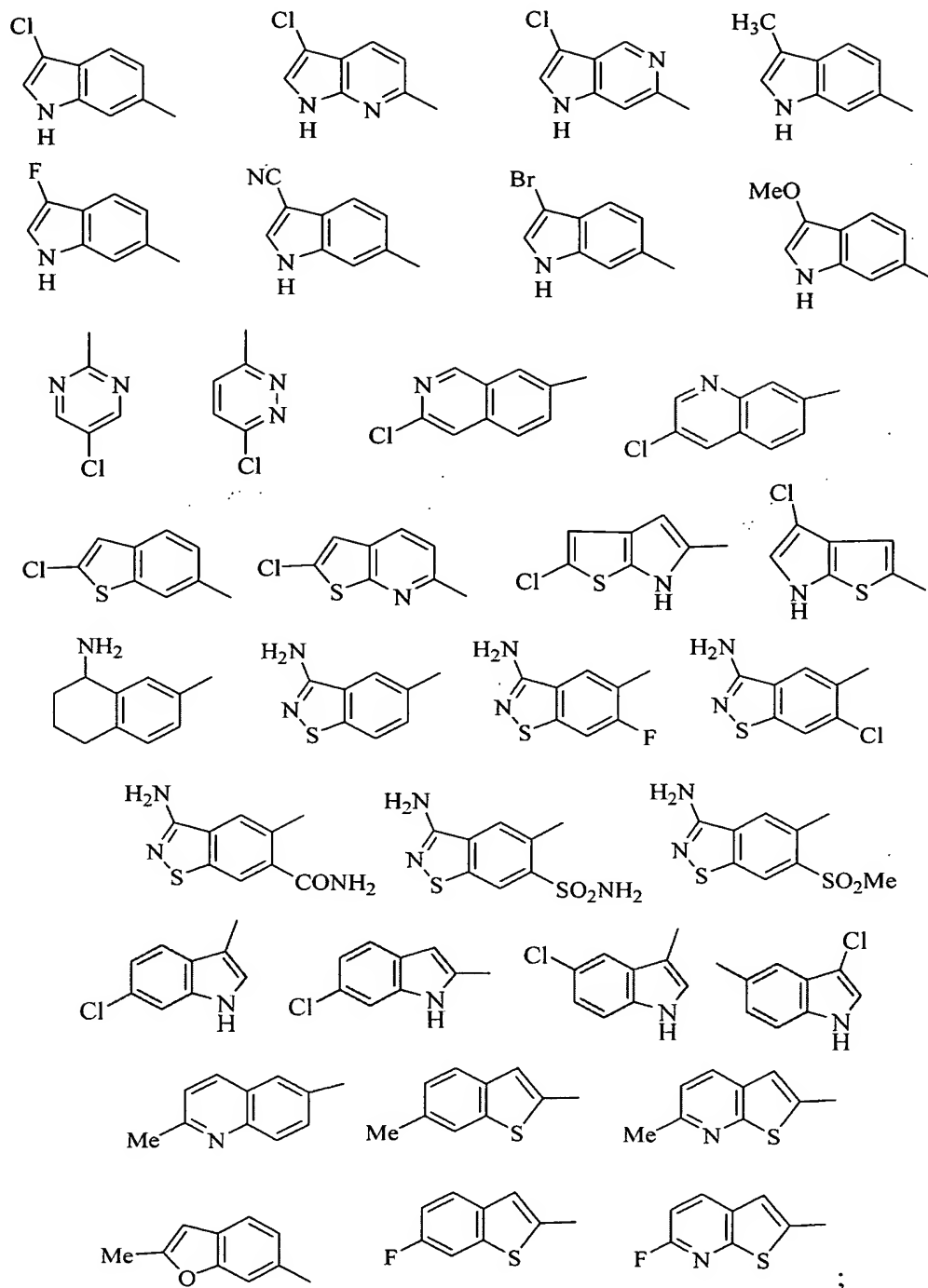
4-bromo-pyrid-2-yl; 4-methyl-pyrid-2-yl; 5-fluoro-thien-2-yl; 5-bromo-thien-2-yl;

- 5-methyl-thien-2-yl; 2-amido-4-methoxy-phenyl; 2-amido-phenyl;
 2-aminomethyl-3-fluoro-phenyl; 2-aminomethyl-4-fluoro-phenyl;
 2-aminomethyl-4-methoxy-phenyl; 2-aminomethyl-5-fluoro-phenyl;
 2-aminomethyl-5-methoxy-phenyl; 2-aminomethyl-6-fluoro-phenyl;
 5 2-aminomethyl-phenyl; 2-amino-pyrid-4-yl; 2-aminosulfonyl-4-methoxy-phenyl;
 2-aminosulfonyl-phenyl; 2-methylsulfonyl-phenyl;
 3-(N,N-dimethylamino)-4-chloro-phenyl; 3-(N,N-dimethylamino)-phenyl;
 3-(N-methylamino)-4-chloro-phenyl; 3-(N-methylamino)-phenyl; 3-amido-phenyl;
 3-amino-4-chloro-phenyl; 3-aminomethyl-phenyl; 3-amino-phenyl; 3-chloro-phenyl;
 10 4-(N,N-dimethylamino)-5-chloro-thien-2-yl; 4-(N-methylamino)-5-chloro-thien-2-yl;
 4-amino-5-chloro-thien-2-yl; 4-chloro-phenyl; 4-methoxy-2-methylsulfonyl-phenyl;
 4-methoxy-phenyl; 2-methoxy-pyrid-5-yl;
 5-(N,N-dimethylamino)-4-chloro-thien-2-yl; 5-(N-methylamino)-4-chloro-thien-2-yl;
 5-amino-4-chloro-thien-2-yl; 5-chloro-pyrid-2-yl; 5-chloro-thien-2-yl;
 15 6-amino-5-chloro-pyrid-2-yl; 6-amino-pyrid-2-yl; 2-cyano-4-chloro-phenyl;
 2-methoxy-4-chloro-phenyl; 2-fluoro-4-chloro-phenyl; phenyl; 4-ethyl-phenyl;
 3-chloro-4-methyl-phenyl; 4-fluoro-phenyl; 3-fluoro-4-chloro-phenyl;
 3-methyl-4-chloro-phenyl; 3-fluoro-4-methyl-phenyl; 3,4-dimethyl-phenyl;
 3-chloro-4-fluoro-phenyl; 3-methyl-4-fluoro-phenyl; 4-methylsulfonyl-phenyl;
 20 2-chlorothiazol-5-yl; 5-chlorothiazol-2-yl;





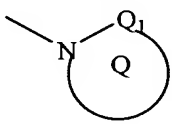




5

A is selected from one of the following carbocyclic and heterocyclic groups which are substituted with 0-2 R⁴; cyclohexyl, phenyl, piperidinyl, piperazinyl, pyridyl, pyrimidyl, furanyl, morpholinyl, thienyl, pyrrolyl, pyrrolidinyl, oxazolyl, isoxazolyl, thiazolyl, isothiazolyl, pyrazolyl, imidazolyl, 1,2,3-oxadiazolyl,

- 1,2,4-oxadiazolyl, 1,2,5-oxadiazolyl, 1,3,4-oxadiazolyl, 1,2,3-thiadiazolyl, 1,2,4-thiadiazolyl, 1,2,5-thiadiazolyl, 1,3,4-thiadiazolyl, 1,2,3-triazolyl, 1,2,4-triazolyl, 1,2,5-triazolyl, 1,3,4-triazolyl, benzofuranyl, benzothiofuranyl, indolinyl, indolyl, benzimidazolyl, benzoxazolyl, benzthiazolyl, indazolyl, 5 benzisoxazolyl, benzisothiazolyl, and isoindazolyl;

B is ; provided that Z and B are attached to different atoms on A; Q₁ is selected from C=O and SO₂;

- ring Q is a 5-7 membered monocyclic or tricyclic ring consisting of, in addition to the N-Q₁ group shown, carbon atoms and 0-1 heteroatoms selected from 10 NR^{4c}, O, and S(O)_p, wherein: 0-2 double bonds are present within the ring and the ring is substituted with 0-2 R^{4a};

- alternatively, ring Q is a 5-7 membered ring to which another ring is fused, wherein: the 5-7 membered ring consists of, in addition to the N-Q₁ group shown, carbon atoms and 0-1 heteroatoms selected from NR^{4c}, O, and S(O)_p and 0-1 double 15 bonds are present within the ring; the fusion ring is phenyl;

ring Q, which includes the 5-7 membered ring and the fusion ring, is substituted with 0-2 R^{4a};

- G₁ is selected from (CR^{3R3a})₁₋₃, CR³=CR³, (CR^{3R3a})_uC(O)(CR^{3R3a})_w, (CR^{3R3a})_uO(CR^{3R3a})_w, (CR^{3R3a})_uNR^{3b}(CR^{3R3a})_w, 20 (CR^{3R3a})_uC(O)NR^{3b}(CR^{3R3a})_w, (CR^{3R3a})_uNR^{3b}C(O)(CR^{3R3a})_w, (CR^{3R3a})_uS(CR^{3R3a})_w, (CR^{3R3a})_uS(O)(CR^{3R3a})_w, (CR^{3R3a})_uS(O)₂(CR^{3R3a})_w, (CR^{3R3a})_uS(O)NR^{3b}(CR^{3R3a})_w, (CR^{3R3a})_uNR^{3b}S(O)₂(CR^{3R3a})_w, (CR^{3R3a})_uS(O)₂NR^{3b}(CR^{3R3a})_w, (CR^{3R3a})_uC(O)NR^{3b}S(O)₂(CR^{3R3a})_w, (CR^{3R3a})_uNR^{3b}C(O)(CR^{3R3a})_uC(O)NR^{3b}(CR^{3R3a})_w, 25 (CR^{3R3a})_uNR^{3b}(CR^{3R3a})_uC(O)NR^{3b}(CR^{3R3a})_w, (CR^{3R3a})_uNR^{3b}C(O)(CR^{3R3a})_uC(S)NR^{3b}(CR^{3R3a})_w, and (CR^{3R3a})_uNR^{3b}C(S)(CR^{3R3a})_uC(O)NR^{3b}(CR^{3R3a})_w, wherein u+w or u+u+w total 0,

1, or 2 and the right side of G_1 is attached to ring G, provided that G_1 does not form a N-S, NCH_2N , NCH_2O , or NCH_2S bond with either group to which it is attached;

Z is selected from $(CR^3R^{3a})_{1-3}$, $(CR^3R^{3a})_u C(O)(CR^3R^{3a})_w$,

$(CR^3R^{3a})_u O(CR^3R^{3a})_w$, $(CR^3R^{3a})_u NR^{3b}(CR^3R^{3a})_w$,

5 $(CR^3R^{3a})_u C(O)NR^{3b}(CR^3R^{3a})_w$, $(CR^3R^{3a})_u NR^{3b}C(O)(CR^3R^{3a})_w$,

$(CR^3R^{3a})_u NR^{3b}C(O)(CR^3R^{3a})_u C(O)NR^{3b}(CR^3R^{3a})_w$, $(CR^3R^{3a})_u S(CR^3R^{3a})_w$,

$(CR^3R^{3a})_u S(O)(CR^3R^{3a})_w$, $(CR^3R^{3a})_u S(O)_2(CR^3R^{3a})_w$,

$(CR^3R^{3a})_u S(O)NR^{3b}(CR^3R^{3a})_w$, $(CR^3R^{3a})_u NR^{3b}S(O)_2(CR^3R^{3a})_w$,

$(CR^3R^{3a})_u S(O)_2NR^{3b}(CR^3R^{3a})_w$, and $(CR^3R^{3a})_u C(O)NR^{3b}S(O)_2(CR^3R^{3a})_w$,

10 wherein $u+w$ or $u+u+w$ total 0, 1, or 2 and the right side of Z is attached to A, provided that G_1 does not form a N-S, NCH_2N , NCH_2O , or NCH_2S bond with either group to which it is attached;

R^{1a} is selected from H, $-(CH_2)_r-R^{1b}$, $-(CH(CH_3))_r-R^{1b}$, $-(C(CH_3)_2)_r-R^{1b}$,

$NHCH_2R^{1c}$, OCH_2R^{1c} , SCH_2R^{1c} , $NH(CH_2)_2(CH_2)_tR^{1b}$, and $O(CH_2)_2(CH_2)_tR^{1b}$,

15 provided that R^{1a} forms other than an N-halo, N-S, or N-CN bond;

alternatively, when two R^{1a} groups are attached to adjacent atoms, together with the atoms to which they are attached they form a 5-7 membered ring consisting of: carbon atoms and 0-2 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, this ring being substituted with 0-2 R^{4b} and having 0-3 ring double bonds;

20 R^{1b} is selected from H, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, F, Cl, Br, I, -CN, -CHO, CF_3 , OR^2 , NR^2R^{2a} , $C(O)R^{2b}$, CO_2R^{2b} , $OC(O)R^2$, CO_2R^{2a} , $S(O)_pR^{2b}$, $NR^2(CH_2)_rOR^2$, $NR^2C(O)R^{2b}$, $NR^2C(O)NHR^2$, $NR^2C(O)_2R^{2a}$, $OC(O)NR^2R^{2a}$, $C(O)NR^2R^{2a}$, $C(O)NR^2R^{2b}$, $C(S)NR^2R^{2a}$, $C(O)NR^2(CH_2)_rOR^2$, $SO_2NR^2R^{2a}$, $NR^2SO_2R^2$, C_{3-6} carbocycle substituted with 0-2 R^{4b} , and 5-6 membered heterocycle
25 consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{4b} , provided that R^{1b} forms other than an O-O, N-halo, N-S, or N-CN bond;

R^{1c} is selected from H, $CH(CH_2OR^2)_2$, $C(O)R^{2c}$, $C(O)NR^2R^{2a}$, $S(O)R^2$,

$S(O)_2R^2$, and $SO_2NR^2R^{2a}$;

R^2 , at each occurrence, is selected from H, CF_3 , CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, benzyl substituted with 0-2 R^{4b} , C_{5-6} carbocycle substituted with 0-2 R^{4b} , a - CH_2 - C_{5-6} carbocyclic group substituted with 0-2 R^{4b} , and 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{4b} ;

R^{2a} , at each occurrence, is selected from H, CF_3 , CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, benzyl substituted with 0-2 R^{4b} , C_{3-6} carbocycle substituted with 0-2 R^{4b} , and 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{4b} ;

alternatively, NR^2R^{2a} forms a 4, 5, or 6 membered saturated, partially saturated, or unsaturated ring substituted with 0-2 R^{4b} and consisting of: carbon atoms, the nitrogen atom to which R^2 and R^{2a} are attached, and 0-1 additional heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^{2b} , at each occurrence, is selected from CF_3 , C_{1-4} alkoxy, C_{1-5} alkyl substituted with 0-3 R^{4b} , benzyl substituted with 0-2 R^{4b} , C_{3-6} carbocycle substituted with 0-2 R^{4b} , and 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{4b} ;

R^{2c} , at each occurrence, is selected from CF_3 , OH, C_{1-4} alkoxy, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, benzyl substituted with 0-2 R^{4b} , C_{5-6} carbocycle substituted with 0-2 R^{4b} , and 5-6 membered heterocycle containing from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{4b} ;

R^3 , at each occurrence, is selected from H, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, benzyl, and phenyl;

R^{3a} , at each occurrence, is selected from H, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, benzyl, and phenyl;

alternatively, NR^3R^{3a} forms a 5 or 6 membered saturated, partially unsaturated, or unsaturated ring consisting of: carbon atoms and the nitrogen atom to which R^3 and R^{3a} are attached;

R^{3c} , at each occurrence, is selected from CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, benzyl, and phenyl;

R^{3d} , at each occurrence, is selected from H, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, CH_2 -phenyl, CH_2CH_2 -phenyl, and $C(=O)R^{3c}$;

R^4 , at each occurrence, is selected from H, $=O$, OR^2 , CH_2OR^2 , $(CH_2)_2OR^2$, F, Cl, Br, I, C_{1-4} alkyl, $-CN$, NO_2 , NR^2R^{2a} , $CH_2NR^2R^{2a}$, $(CH_2)_2NR^2R^{2a}$, $C(O)R^{2c}$, $NR^2C(O)R^{2b}$, $C(O)NR^2R^{2a}$, $SO_2NR^2R^{2a}$, $S(O)_pR^{5a}$, CF_3 , CF_2CF_3 , 5-6 membered carbocycle substituted with 0-1 R^5 , and a 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-1 R^5 ;

R^{4a} , at each occurrence, is selected from H, $=O$, OR^2 , CH_2OR^2 , F, CH_2F , Br, CH_2Br , Cl, CH_2Cl , C_{1-4} alkyl, $-CN$, $-CH_2CN$, NO_2 , CH_2NO_2 , NR^2R^{2a} , $CH_2NR^2R^{2a}$, $C(O)R^{2c}$, $CH_2C(O)R^{2c}$, $NR^2C(O)R^{2b}$, $(CH_2)_rC(O)NR^2R^{2a}$, $NR^2C(O)NR^2R^{2a}$, $(CH_2)_rSO_2NR^2R^{2a}$, $NR^2SO_2NR^2R^{2a}$, $NR^2SO_2-C_{1-4}$ alkyl, $NR^2SO_2R^5$, $(CH_2)_rS(O)_pR^{5a}$, CH_2CF_3 , CF_3 , CH_2 -5-6 membered carbocycle substituted with 0-1 R^5 , 5-6 membered carbocycle substituted with 0-1 R^5 , and a CH_2 -5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-1 R^5 , and 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-1 R^5 ;

R^{4b} , at each occurrence, is selected from H, $=O$, OR^3 , CH_2OR^3 , F, Cl, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, $-CN$, NO_2 , NR^3R^{3a} , $CH_2NR^3R^{3a}$, $C(O)R^3$, $CH_2C(O)R^3$, $C(O)OR^{3c}$, $CH_2C(O)OR^{3c}$, $NR^3C(O)R^{3a}$, $CH_2NR^3C(O)R^{3a}$,

$C(O)NR^3R^{3a}$, $CH_2C(O)NR^3R^{3a}$, $NR^3C(O)NR^3R^{3a}$, $CH_2NR^3C(O)NR^3R^{3a}$,
 $C(=NR^3)NR^3R^{3a}$, $CH_2C(=NR^3)NR^3R^{3a}$, $NR^3C(=NR^3)NR^3R^{3a}$,
 $CH_2NR^3C(=NR^3)NR^3R^{3a}$, $SO_2NR^3R^{3a}$, $CH_2SO_2NR^3R^{3a}$, $NR^3SO_2NR^3R^{3a}$,
 $CH_2NR^3SO_2NR^3R^{3a}$, $NR^3SO_2-C_{1-4}$ alkyl, $CH_2NR^3SO_2-C_{1-4}$ alkyl, $NR^3SO_2CF_3$,
5 $CH_2NR^3SO_2CF_3$, NR^3SO_2 -phenyl, $CH_2NR^3SO_2$ -phenyl, $S(O)_pCF_3$, $CH_2S(O)_pCF_3$,
 $S(O)_p-C_{1-4}$ alkyl, $CH_2S(O)_p-C_{1-4}$ alkyl, $S(O)_p$ -phenyl, $CH_2S(O)_p$ -phenyl, CF_3 , and
 CH_2CF_3 ;

R^{4c} , at each occurrence, is selected from H, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$,
 $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$,
10 CH_2OR^2 , CH_2F , CH_2Br , CH_2Cl , CH_2CN , CH_2NO_2 , $CH_2NR^2R^{2a}$, $C(O)R^{2c}$,
 $CH_2C(O)R^{2c}$, $CH_2NR^2C(O)R^{2b}$, $C(O)NR^2R^{2a}$, $CH_2C(O)NR^2R^{2a}$,
 $CH_2NR^2C(O)NR^2R^{2a}$, $SO_2NR^2R^{2a}$, $CH_2SO_2NR^2R^{2a}$, $CH_2NR^2SO_2NR^2R^{2a}$,
 $CH_2NR^2SO_2-C_{1-4}$ alkyl, $C(O)NHSO_2-C_{1-4}$ alkyl, $CH_2C(O)NHSO_2-C_{1-4}$ alkyl,
 $CH_2NR^2SO_2R^5$, $S(O)_pR^{5a}$, $CH_2S(O)_pR^{5a}$, CF_3 , CH_2CF_3 , 5-6 membered carbocycle
15 substituted with 0-1 R^5 , CH_2 -5-6 membered carbocycle substituted with 0-1 R^5 , 5-6
membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected
from the group consisting of N, O, and $S(O)_p$, and substituted with 0-1 R^5 , and a
 CH_2 -5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms
selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-1 R^5 ;

R^5 , at each occurrence, is selected from H, $=O$, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$,
 $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, OR^3 ,
 CH_2OR^3 , F, Cl, -CN, NO_2 , NR^3R^{3a} , $CH_2NR^3R^{3a}$, $C(O)R^3$, $CH_2C(O)R^3$, $C(O)OR^{3c}$,
 $CH_2C(O)OR^{3c}$, $NR^3C(O)R^{3a}$, $C(O)NR^3R^{3a}$, $NR^3C(O)NR^3R^{3a}$, $CH(=NOR^{3d})$,
 $C(=NR^3)NR^3R^{3a}$, $NR^3C(=NR^3)NR^3R^{3a}$, $SO_2NR^3R^{3a}$, $NR^3SO_2NR^3R^{3a}$,
25 $NR^3SO_2-C_{1-4}$ alkyl, $NR^3SO_2CF_3$, NR^3SO_2 -phenyl, $S(O)_pCF_3$, $S(O)_p-C_{1-4}$ alkyl,
 $S(O)_p$ -phenyl, CF_3 , phenyl substituted with 0-2 R^6 , naphthyl substituted with 0-2 R^6 ,
and benzyl substituted with 0-2 R^6 ;

R^{5a} , at each occurrence, is selected from C_{1-6} alkyl, OR^3 , CH_2OR^3 , NR^3R^{3a} ,
 $CH_2NR^3R^{3a}$, $C(O)R^3$, $CH_2C(O)R^3$, $C(O)OR^{3c}$, $CH_2C(O)OR^{3c}$, $NR^3C(O)R^{3a}$,

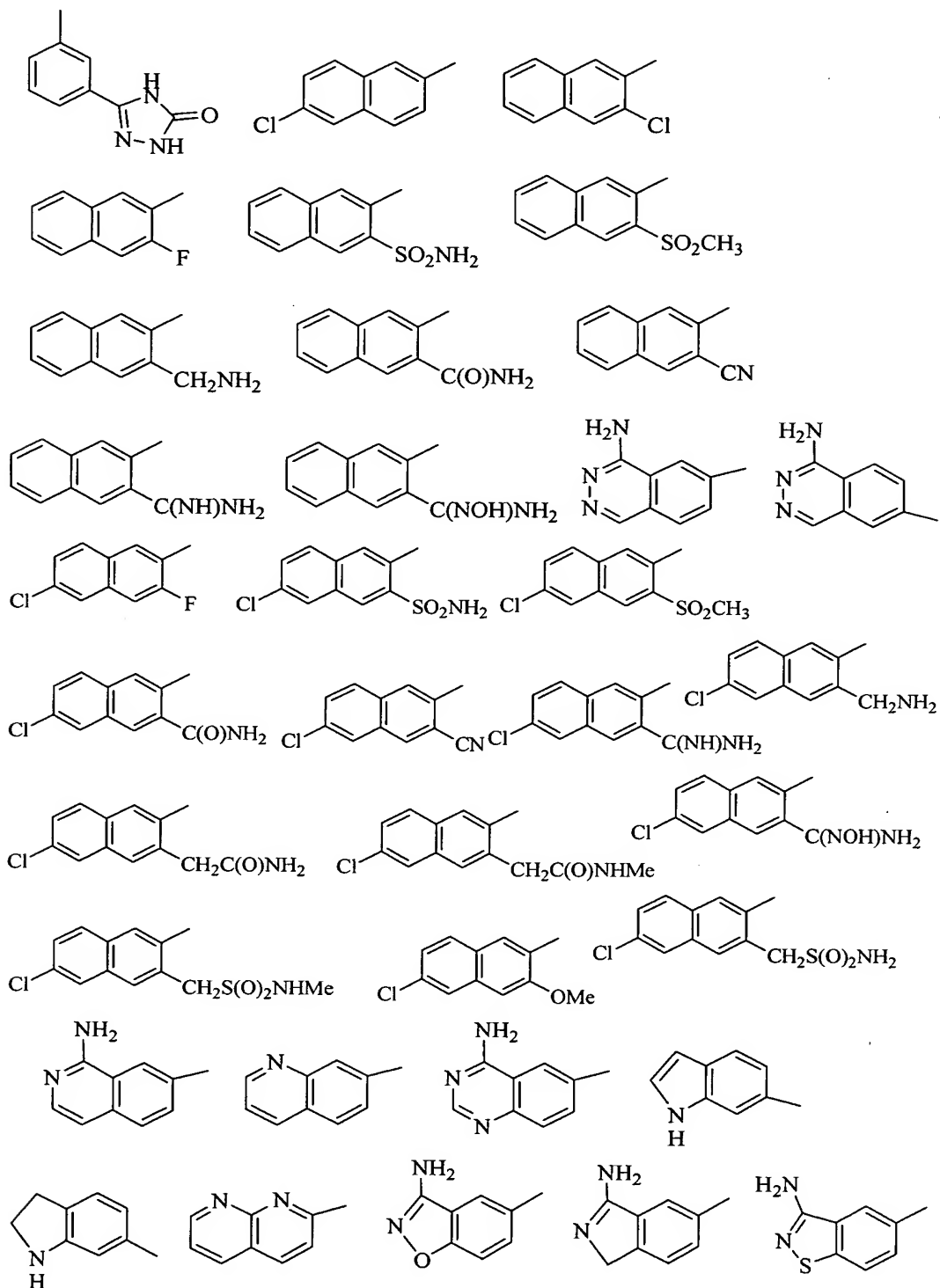
$\text{CH}_2\text{NR}^3\text{C}(\text{O})\text{R}^{3a}$, $\text{C}(\text{O})\text{NR}^3\text{R}^{3a}$, $\text{CH}_2\text{C}(\text{O})\text{NR}^3\text{R}^{3a}$, CF_3 , CF_2CF_3 , phenyl substituted with 0-2 R^6 , naphthyl substituted with 0-2 R^6 , and benzyl substituted with 0-2 R^6 , provided that R^{5a} does not form a S-N or $\text{S}(\text{O})_p\text{-C}(\text{O})$ bond; and

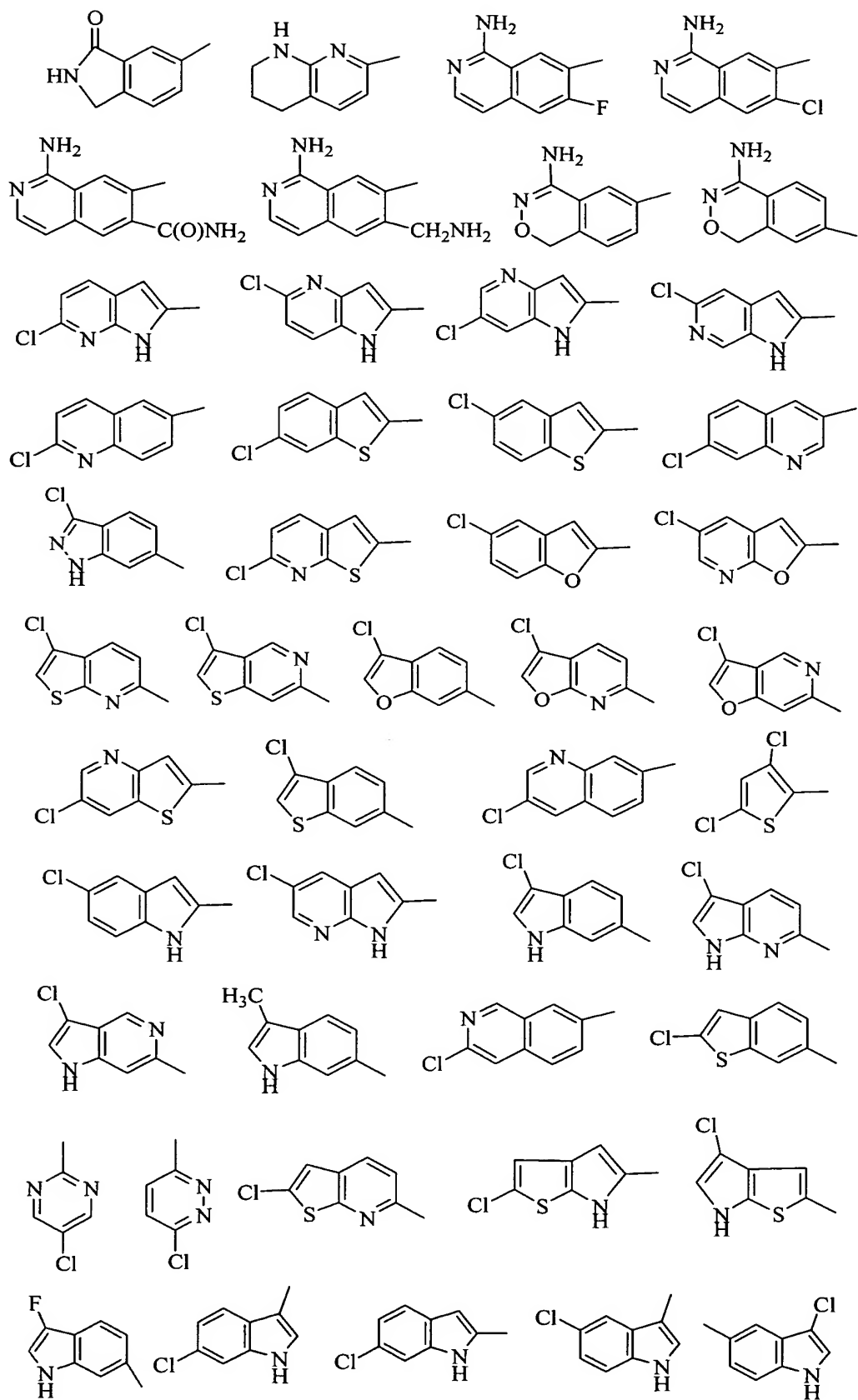
R^6 , at each occurrence, is selected from H, OH, OR^2 , F, Cl, CH_3 , CH_2CH_3 ,
 5 $\text{CH}_2\text{CH}_2\text{CH}_3$, $\text{CH}(\text{CH}_3)_2$, $\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_3$, $\text{CH}_2\text{CH}(\text{CH}_3)_2$, $\text{CH}(\text{CH}_3)\text{CH}_2\text{CH}_3$,
 $\text{C}(\text{CH}_3)_3$, -CN, NO_2 , NR^2R^{2a} , $\text{CH}_2\text{NR}^2\text{R}^{2a}$, $\text{C}(\text{O})\text{R}^{2b}$, $\text{CH}_2\text{C}(\text{O})\text{R}^{2b}$, $\text{NR}^2\text{C}(\text{O})\text{R}^{2b}$,
 $\text{NR}^2\text{C}(\text{O})\text{NR}^2\text{R}^{2a}$, $\text{C}(=\text{NH})\text{NH}_2$, $\text{NHC}(=\text{NH})\text{NH}_2$, $\text{SO}_2\text{NR}^2\text{R}^{2a}$, $\text{NR}^2\text{SO}_2\text{NR}^2\text{R}^{2a}$,
 and $\text{NR}^2\text{SO}_2\text{C}_{1-4}$ alkyl.

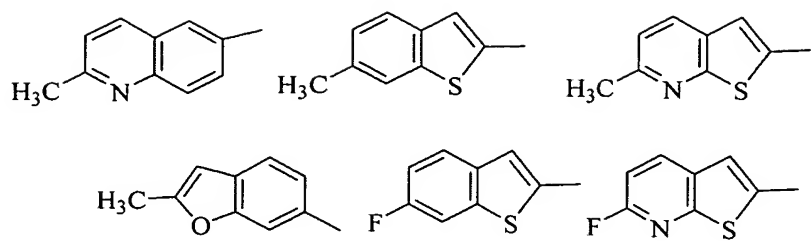
10 3. A compound according to Claim 2, wherein:

G is selected from: phenyl; 4-ethyl-phenyl; 2-aminomethyl-4-chloro-phenyl;
 2-aminosulfonyl-4-chloro-phenyl; 2-amido-4-chloro-phenyl;
 4-chloro-2-methylsulfonyl-phenyl; 2-aminosulfonyl-4-fluoro-phenyl;
 2-amido-4-fluoro-phenyl; 4-fluoro-2-methylsulfonyl-phenyl;
 15 2-aminomethyl-4-bromo-phenyl; 2-aminosulfonyl-4-bromo-phenyl;
 2-amido-4-bromo-phenyl; 4-bromo-2-methylsulfonyl-phenyl;
 2-aminomethyl-4-methyl-phenyl; 2-aminosulfonyl-4-methyl-phenyl;
 2-amido-4-methyl-phenyl; 2-methylsulfonyl-4-methyl-phenyl; 4-fluoro-pyrid-2-yl;
 4-bromo-pyrid-2-yl; 4-methyl-pyrid-2-yl; 5-fluoro-thien-2-yl; 5-bromo-thien-2-yl;
 20 5-methyl-thien-2-yl; 2-amido-4-methoxy-phenyl; 2-amido-phenyl;
 2-aminomethyl-3-fluoro-phenyl; 2-aminomethyl-4-fluoro-phenyl;
 2-aminomethyl-5-fluoro-phenyl; 2-aminomethyl-6-fluoro-phenyl;
 2-aminomethyl-phenyl; 2-amino-pyrid-4-yl; 2-aminosulfonyl-4-methoxy-phenyl;
 2-aminosulfonyl-phenyl; 3-amido-phenyl; 3-amino-4-chloro-phenyl;
 25 3-aminomethyl-phenyl; 3-chloro-phenyl; 4-chloro-phenyl; 4-methoxy-phenyl;
 2-methoxy-pyrid-5-yl; 5-chloro-pyrid-2-yl; 5-chloro-thien-2-yl;
 6-amino-5-chloro-pyrid-2-yl; 6-amino-pyrid-2-yl; 2-cyano-4-chloro-phenyl;
 2-methoxy-4-chloro-phenyl; 2-fluoro-4-chloro-phenyl; 3-chloro-4-methyl-phenyl;
 4-fluoro-phenyl; 3-fluoro-4-chloro-phenyl; 3-methyl-4-chloro-phenyl;
 30 3-fluoro-4-methyl-phenyl; 3,4-dimethyl-phenyl; 3-chloro-4-fluoro-phenyl;

3-methyl-4-fluoro-phenyl; 4-methylsulfanyl-phenyl; 2-chlorothiazol-5-yl;
5-chlorothiazol-2-yl;

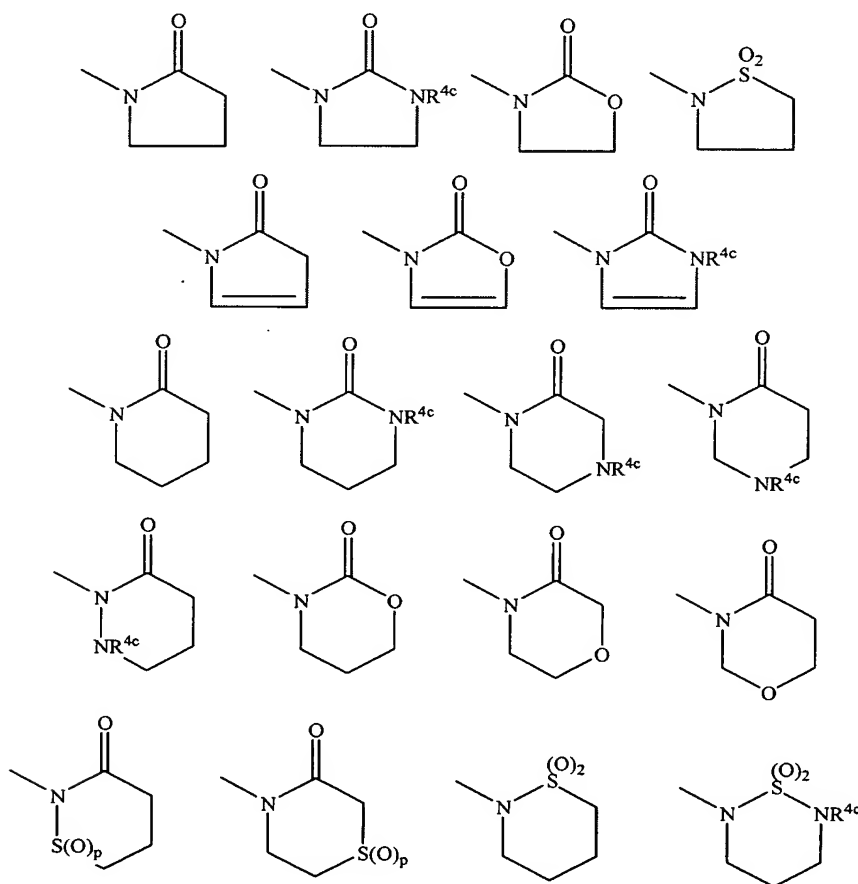




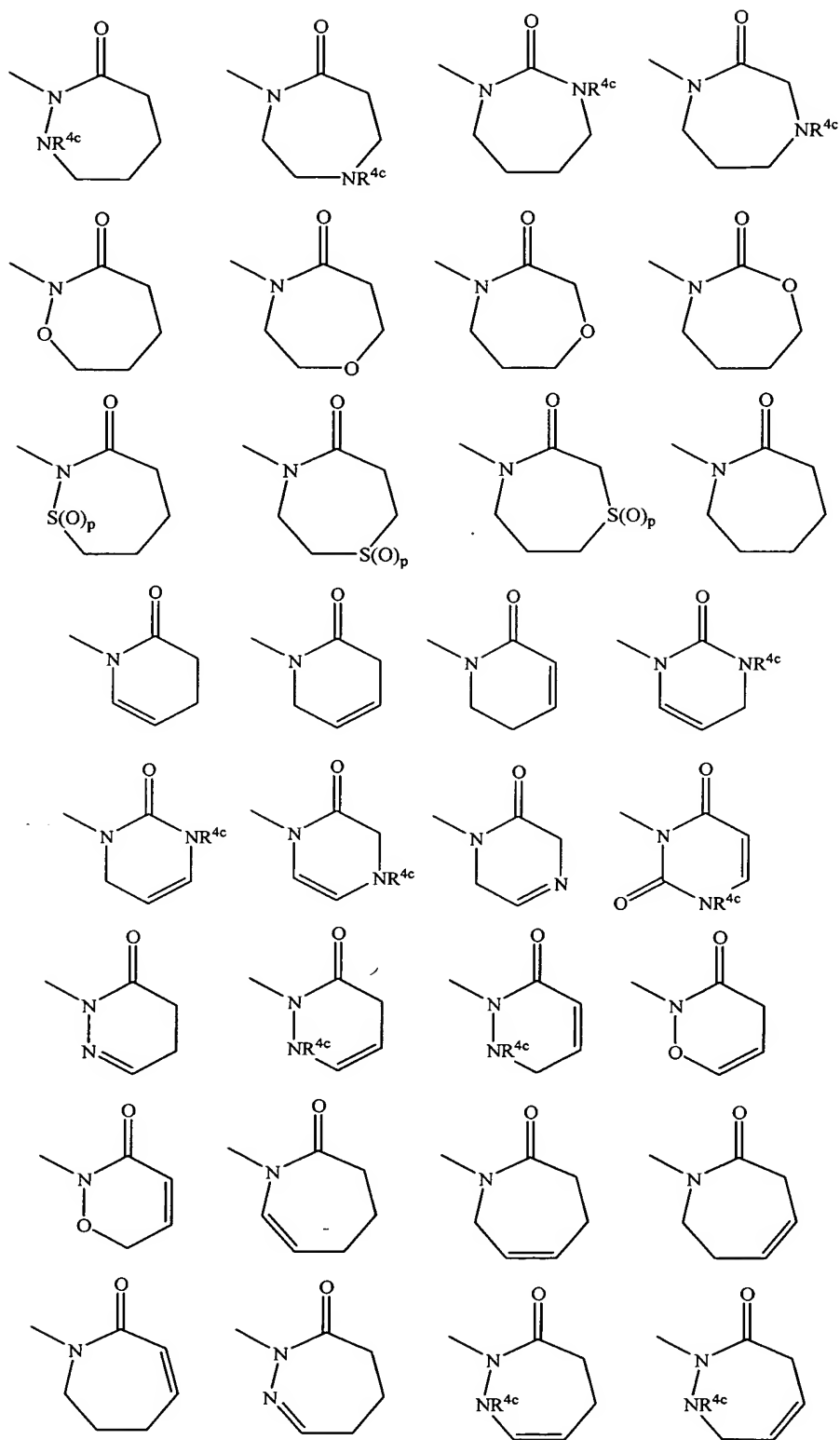


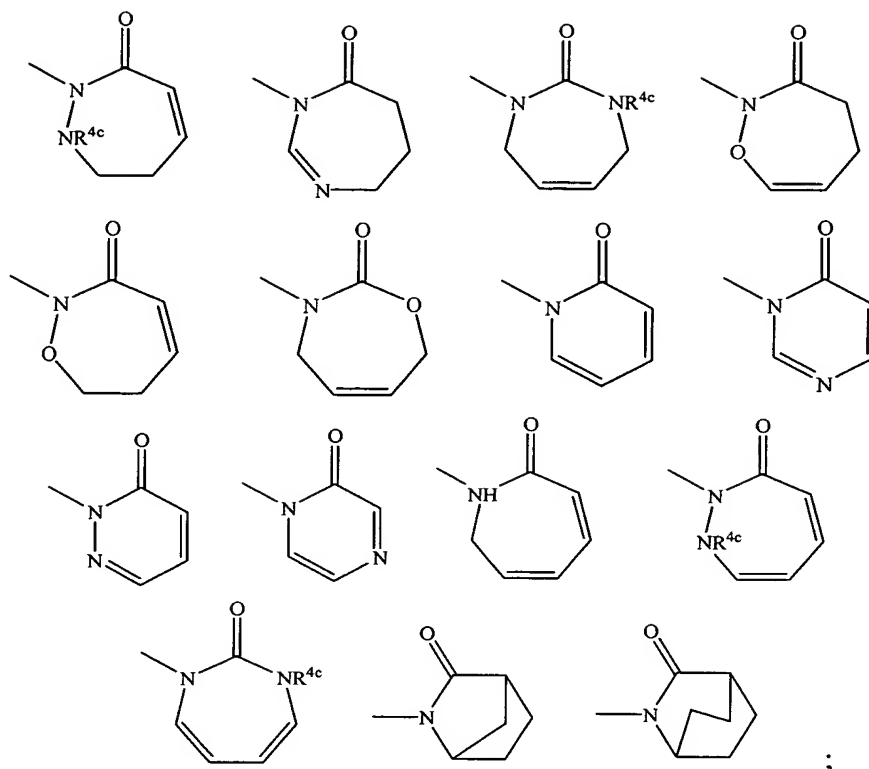
A is selected from the group: cyclohexyl, piperidinyl, indolinyl, phenyl,
 2-pyridyl, 3-pyridyl, 2-pyrimidyl, 2-chloro-phenyl, 3-chloro-phenyl, 2-fluoro-phenyl,
 5 3-fluoro-phenyl, 2-methylphenyl, 3-methylphenyl, 2-aminophenyl, 3-aminophenyl,
 2-methoxyphenyl, and 3-methoxyphenyl;

B is attached to a different atom on A than M, is substituted with 0-2 R^{4a} , and
 is selected from the group:



10





G_1 is selected from CH_2 , CH_2CH_2 , $CH=CH$, CH_2O , OCH_2 , $C(O)$, NH ,
 5 CH_2NH , $NHCH_2$, $CH_2C(O)$, $C(O)CH_2$, $C(O)NH$, $NHC(O)$, $NHC(O)NH$,
 $NHC(O)CH_2C(O)NH$, $C(O)NHS(O)_2$, CH_2S , SCH_2 , $CH_2S(O)$, $S(O)_2$, $CH_2S(O)_2$,
 $S(O)_2(CH_2)$, SO_2NH , $NHSO_2$, $NHCH_2C(O)NH$, $NHC(O)C(O)NH$, $NHC(O)C(S)NH$,
 and $NHC(S)C(O)NH$ and the right side of G_1 is attached to ring G , provided that Z
 does not form a $N-S$, NCH_2N , NCH_2O , or NCH_2S bond with either group to which it
 10 is attached;

Z is selected from CH_2 , CH_2CH_2 , CH_2O , OCH_2 , $C(O)$, NH , CH_2NH ,
 $NHCH_2$, $CH_2C(O)$, $C(O)CH_2$, $C(O)NH$, $NHC(O)$, $NHC(O)NH$,
 $NHC(O)CH_2C(O)NH$, $C(O)NHS(O)_2$, CH_2S , SCH_2 , $CH_2S(O)$, $S(O)_2$, $CH_2S(O)_2$,
 $S(O)_2(CH_2)$, SO_2NH , and $NHSO_2$ and the right side of Z is attached to A , provided
 15 that Z does not form a $N-S$, NCH_2N , NCH_2O , or NCH_2S bond with either group to
 which it is attached;

R^{1a} is selected from H , R^{1b} , $CH(CH_3)R^{1b}$, $C(CH_3)_2R^{1b}$, CH_2R^{1b} , and
 $CH_2CH_2R^{1b}$, provided that R^{1a} forms other than an N -halo, $N-S$, or $N-CN$ bond;

alternatively, when two R^{1a} groups are attached to adjacent atoms, together with the atoms to which they are attached they form a 5-6 membered ring consisting of: carbon atoms and 0-2 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, this ring being substituted with 0-2 R^{4b} and having 0-3 ring double bonds;

- 5 R^{1b} is selected from H, CH_3 , CH_2CH_3 , F, Cl, Br, -CN, -CHO, CF_3 , OR^2 , NR^2R^{2a} , $C(O)R^{2b}$, CO_2R^{2b} , $OC(O)R^2$, CO_2R^{2a} , $S(O)_pR^{2b}$, $NR^2(CH_2)_rOR^2$, $NR^2C(O)R^{2b}$, $C(O)NR^2R^{2a}$, $C(O)NR^2R^{2b}$, $C(S)NR^2R^{2a}$, $SO_2NR^2R^{2a}$, $NR^2SO_2R^2$, C_{3-5} cycloalkyl substituted with 0-2 R^{4b} , phenyl substituted with 0-2 R^{4b} , and 5-6 membered aromatic heterocycle consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{4b} ,
10 provided that R^{1b} forms other than an O-O, N-halo, N-S, or N-CN bond;

- R^2 , at each occurrence, is selected from H, CF_3 , CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, phenyl substituted with 0-2 R^{4b} , a benzyl substituted with 0-2 R^{4b} , and 5-6 membered aromatic heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{4b} ;

- R^{2a} , at each occurrence, is selected from H, CF_3 , CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, $C(CH_3)_3$, C_{3-5} cycloalkyl substituted with 0-1 R^{4b} , benzyl substituted with 0-2 R^{4b} , phenyl substituted with 0-2 R^{4b} , and 5-6 membered aromatic heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{4b} ;

- alternatively, NR^2R^{2a} forms a 4, 5, or 6 membered saturated, partially saturated, or unsaturated ring substituted with 0-2 R^{4b} and consisting of: carbon atoms, the nitrogen atom to which R^2 and R^{2a} are attached, and 0-1 additional heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

- 25 R^{2b} , at each occurrence, is selected from CF_3 , C_{1-4} alkoxy, C_{1-5} alkyl substituted with 0-3 R^{4b} , C_{3-5} cycloalkyl substituted with 0-2 R^{4b} , benzyl substituted with 0-2 R^{4b} , phenyl substituted with 0-2 R^{4b} , and 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{4b} ;

R^{2c} , at each occurrence, is selected from CF_3 , OH, OCH_3 , OCH_2CH_3 , $OCH_2CH_2CH_3$, $OCH(CH_3)_2$, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, benzyl substituted with 0-2 R^{4b} , phenyl substituted with 0-2 R^{4b} , and 5-6 membered aromatic heterocycle containing from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{4b} ;

R^{4a} , at each occurrence, is selected from H, =O, CH_2OR^2 , OR^2 , F, Br, Cl, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, -CN, NO_2 , NR^2R^{2a} , $CH_2NR^2R^{2a}$, $C(O)R^{2c}$, $NR^2C(O)R^{2b}$, $C(O)NR^2R^{2a}$, $NR^2C(O)NR^2R^{2a}$, $SO_2NR^2R^{2a}$, and $-CF_3$;

R^{4b} , at each occurrence, is selected from H, =O, OR^3 , CH_2OR^3 , F, Cl, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, -CN, NO_2 , NR^3R^{3a} , $CH_2NR^3R^{3a}$, $C(O)R^3$, $CH_2C(O)R^3$, $C(O)OR^{3c}$, $CH_2C(O)OR^{3c}$, $NR^3C(O)R^{3a}$, $CH_2NR^3C(O)R^{3a}$, $C(O)NR^3R^{3a}$, $CH_2C(O)NR^3R^{3a}$, $SO_2NR^3R^{3a}$, $CH_2SO_2NR^3R^{3a}$, $NR^3SO_2-C_{1-4}$ alkyl, $CH_2NR^3SO_2-C_{1-4}$ alkyl, NR^3SO_2 -phenyl, $CH_2NR^3SO_2$ -phenyl, $S(O)_pCF_3$, $CH_2S(O)_pCF_3$, $S(O)_p-C_{1-4}$ alkyl, $CH_2S(O)_p-C_{1-4}$ alkyl, $S(O)_p$ -phenyl, $CH_2S(O)_p$ -phenyl, and CF_3 ;

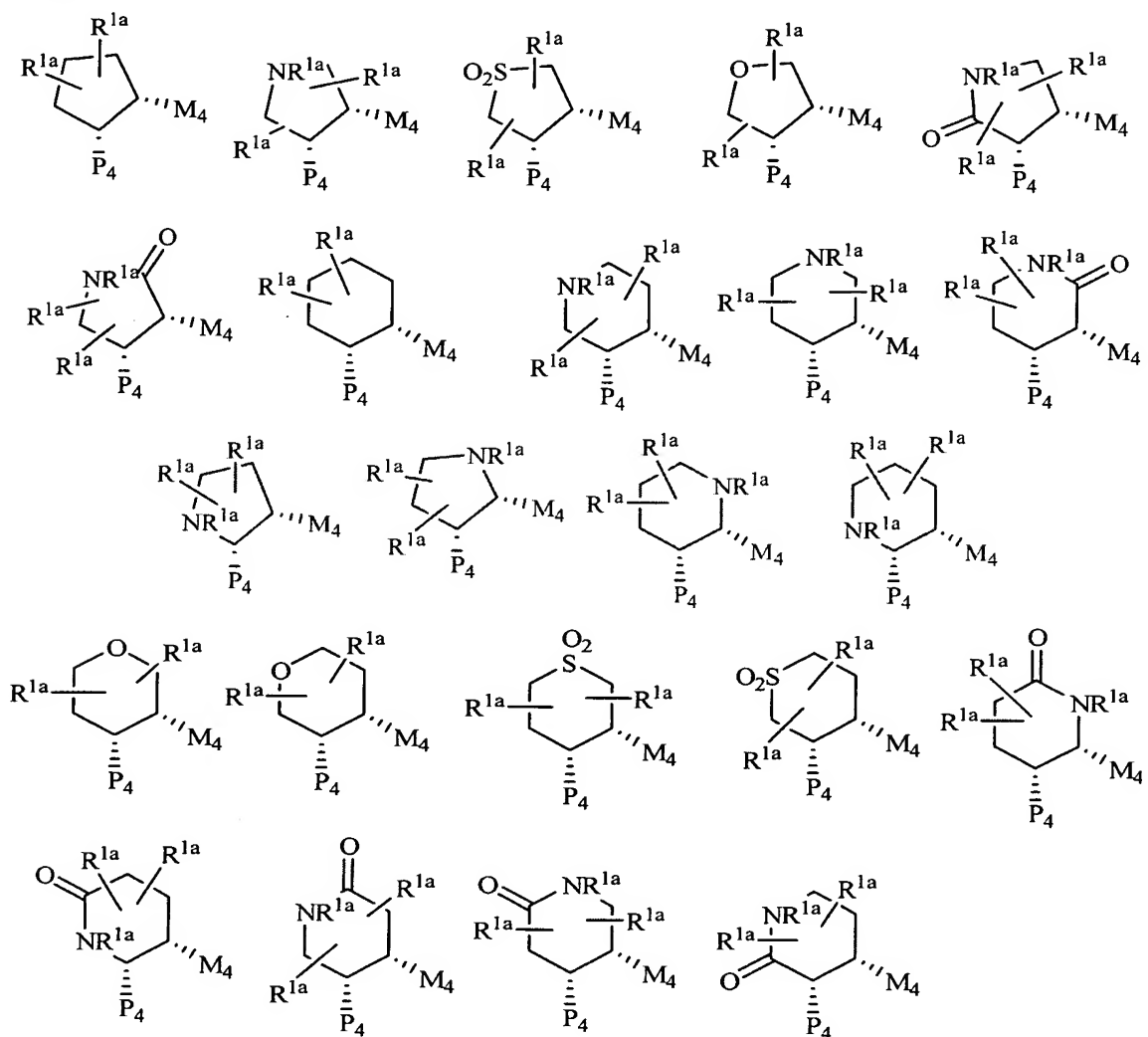
R^{4c} , at each occurrence, is selected from H, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, $CH_2CH_2CH_2CH_3$, $CH_2CH(CH_3)_2$, $CH(CH_3)CH_2CH_3$, $C(CH_3)_3$, CH_2OR^2 , CH_2F , CH_2Br , CH_2Cl , CH_2CN , CH_2NO_2 , $CH_2NR^2R^{2a}$, $C(O)R^{2c}$, $CH_2C(O)R^{2c}$, $CH_2NR^2C(O)R^{2b}$, $C(O)NR^2R^{2a}$, $CH_2C(O)NR^2R^{2a}$, $SO_2NR^2R^{2a}$, $CH_2SO_2NR^2R^{2a}$, $S(O)_pR^{5a}$, $CH_2S(O)_pR^{5a}$, CF_3 , phenyl substituted with 0-1 R^5 , and benzyl substituted with 0-1 R^5 ;

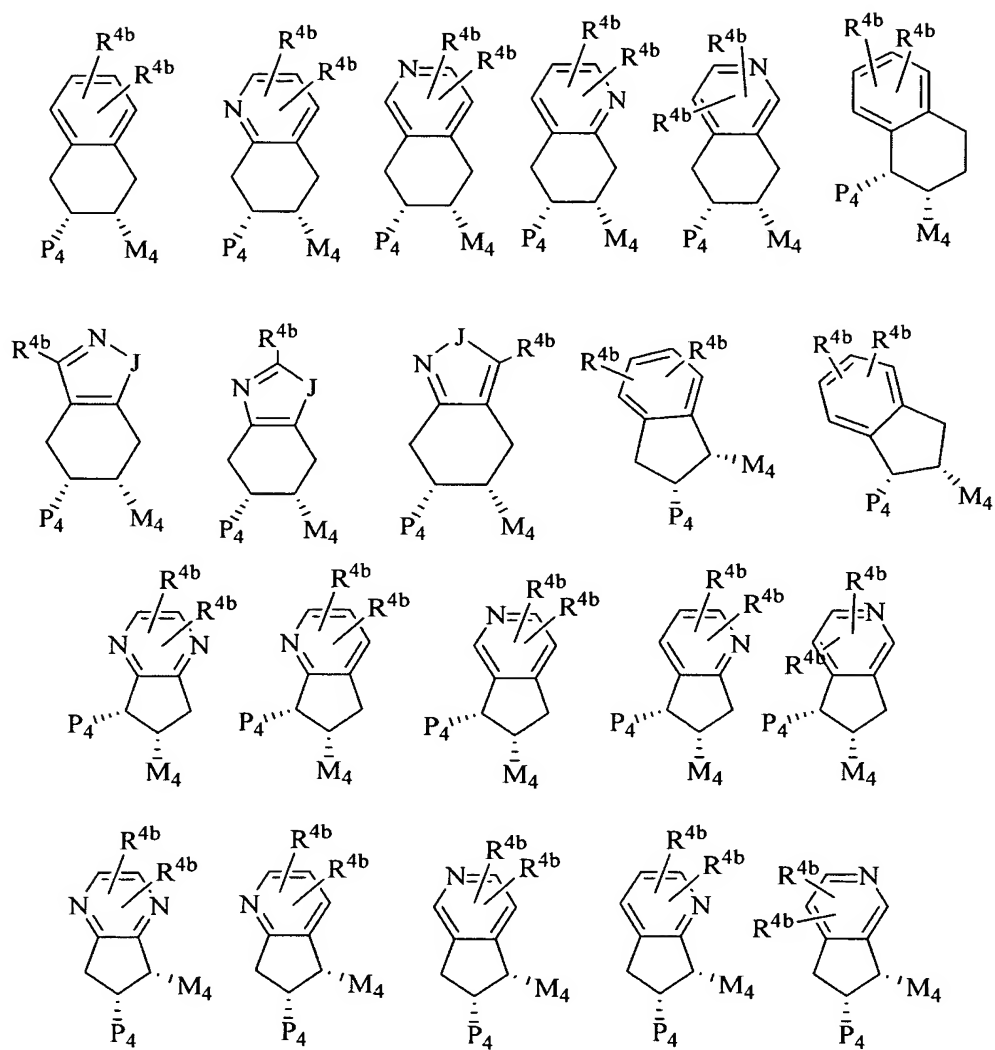
R^5 , at each occurrence, is selected from H, =O, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, OR^3 , CH_2OR^3 , F, Cl, -CN, NO_2 , NR^3R^{3a} , $CH_2NR^3R^{3a}$, $C(O)R^3$, $CH_2C(O)R^3$, $C(O)OR^{3c}$, $CH_2C(O)OR^{3c}$, $NR^3C(O)R^{3a}$, $C(O)NR^3R^{3a}$, $SO_2NR^3R^{3a}$, $NR^3SO_2-C_{1-4}$ alkyl, $NR^3SO_2CF_3$, NR^3SO_2 -phenyl, $S(O)_pCF_3$, $S(O)_p-C_{1-4}$ alkyl, $S(O)_p$ -phenyl, CF_3 , phenyl substituted with 0-2 R^6 , naphthyl substituted with 0-2 R^6 , and benzyl substituted with 0-2 R^6 ;

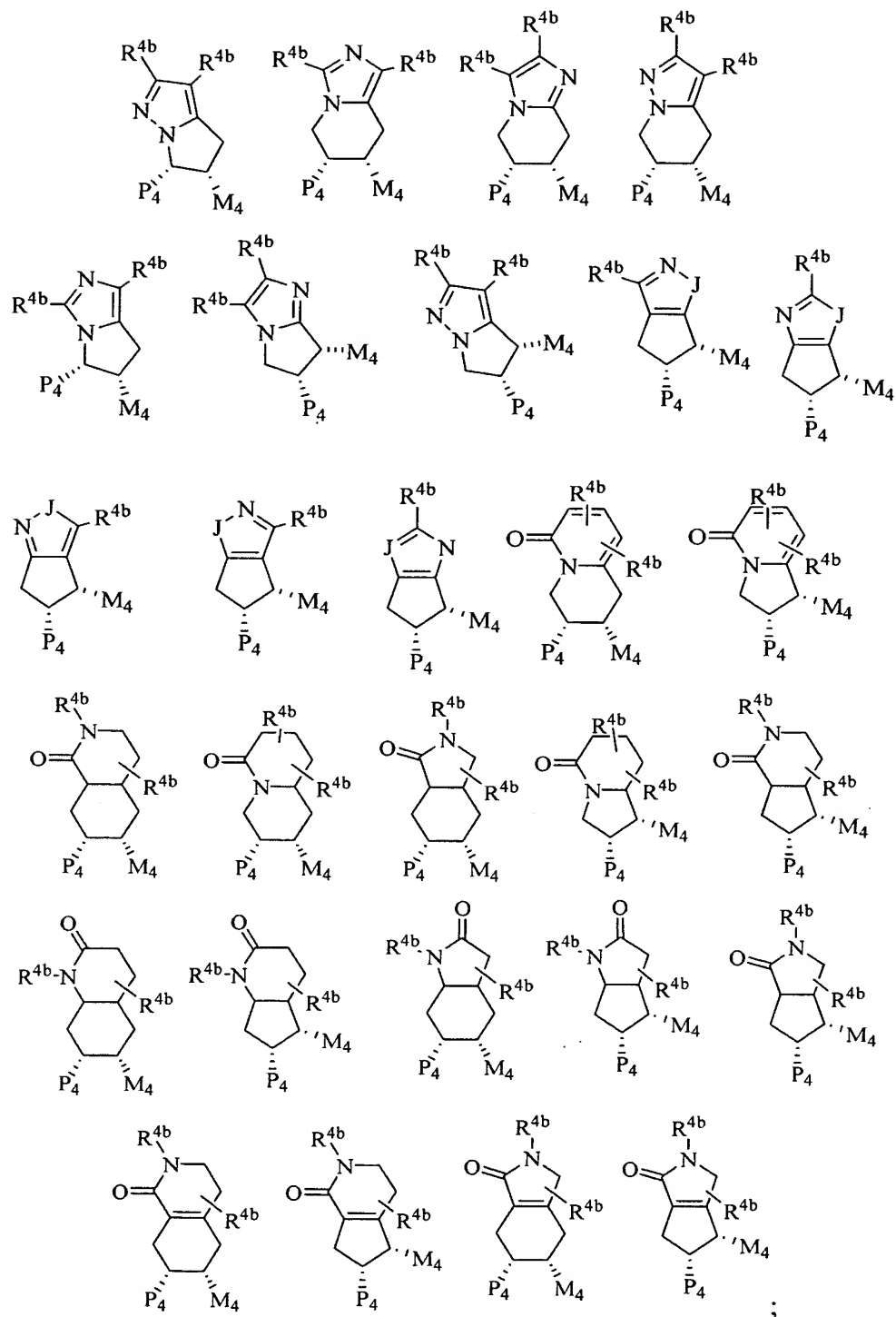
R^{5a} , at each occurrence, is selected from CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, OR^3 , NR^3R^{3a} , $C(O)R^3$, $NR^3C(O)R^{3a}$, $C(O)NR^3R^{3a}$, CF_3 , phenyl substituted with 0-2 R^6 , naphthyl substituted with 0-2 R^6 , and benzyl substituted with 0-2 R^6 , provided that R^{5a} does not form a S-N or $S(O)_p-C(O)$ bond; and

- 5 R^6 , at each occurrence, is selected from H, OH, OR^2 , F, Cl, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, -CN, NO_2 , NR^2R^{2a} , $CH_2NR^2R^{2a}$, $C(O)R^{2b}$, $CH_2C(O)R^{2b}$, $NR^2C(O)R^{2b}$, $SO_2NR^2R^{2a}$, and $NR^2SO_2C_{1-4}$ alkyl.

- 10 4. A compound according to Claim 3, wherein the compound is selected from:



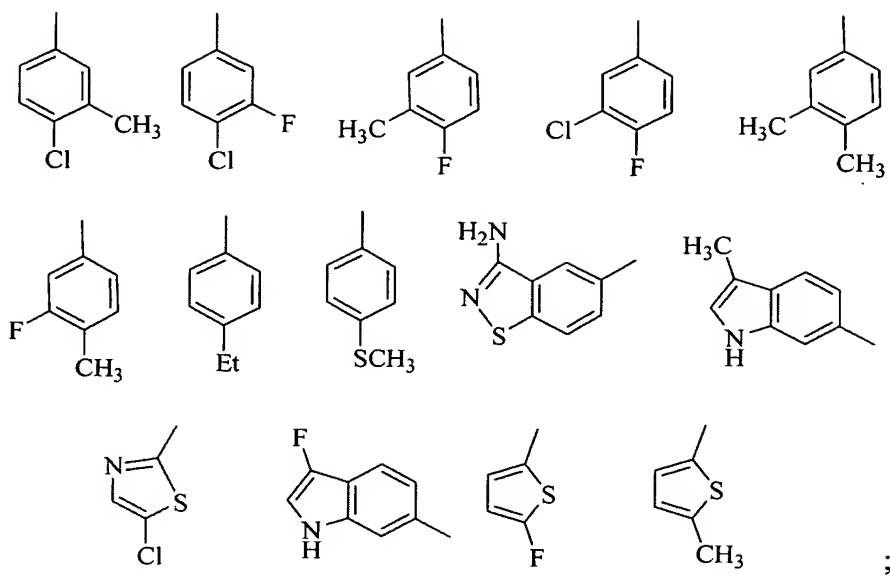




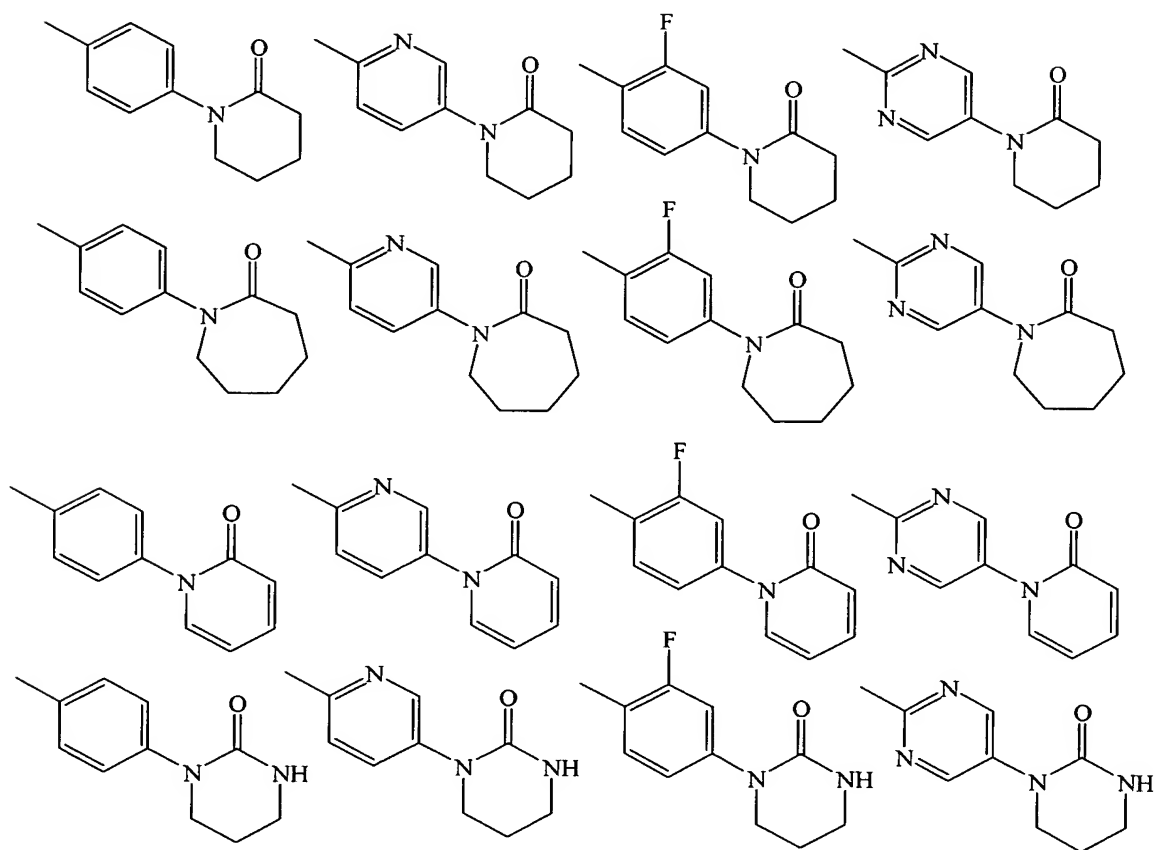
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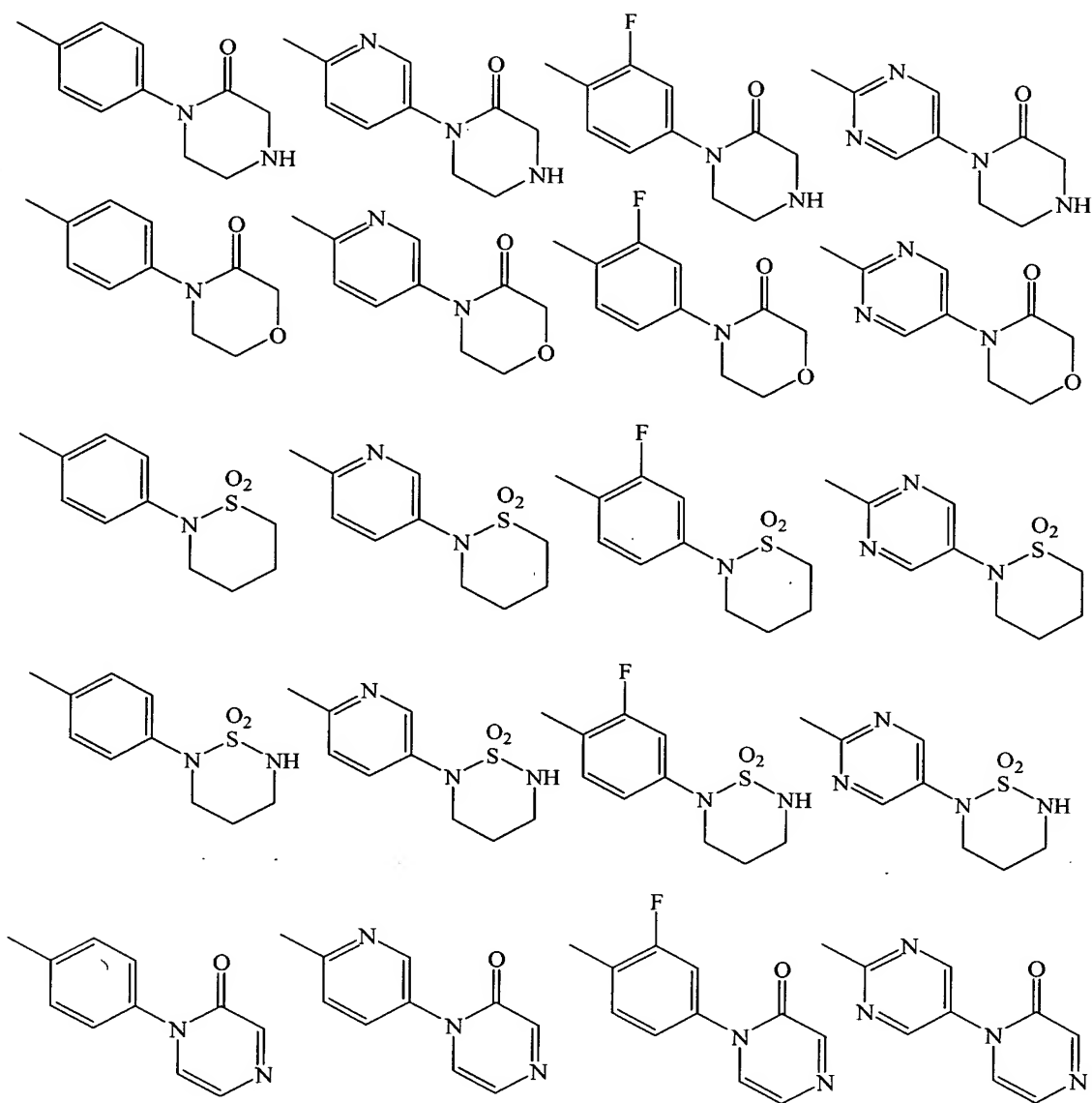
J is selected from O, S, NH, and NR^{1a} ;

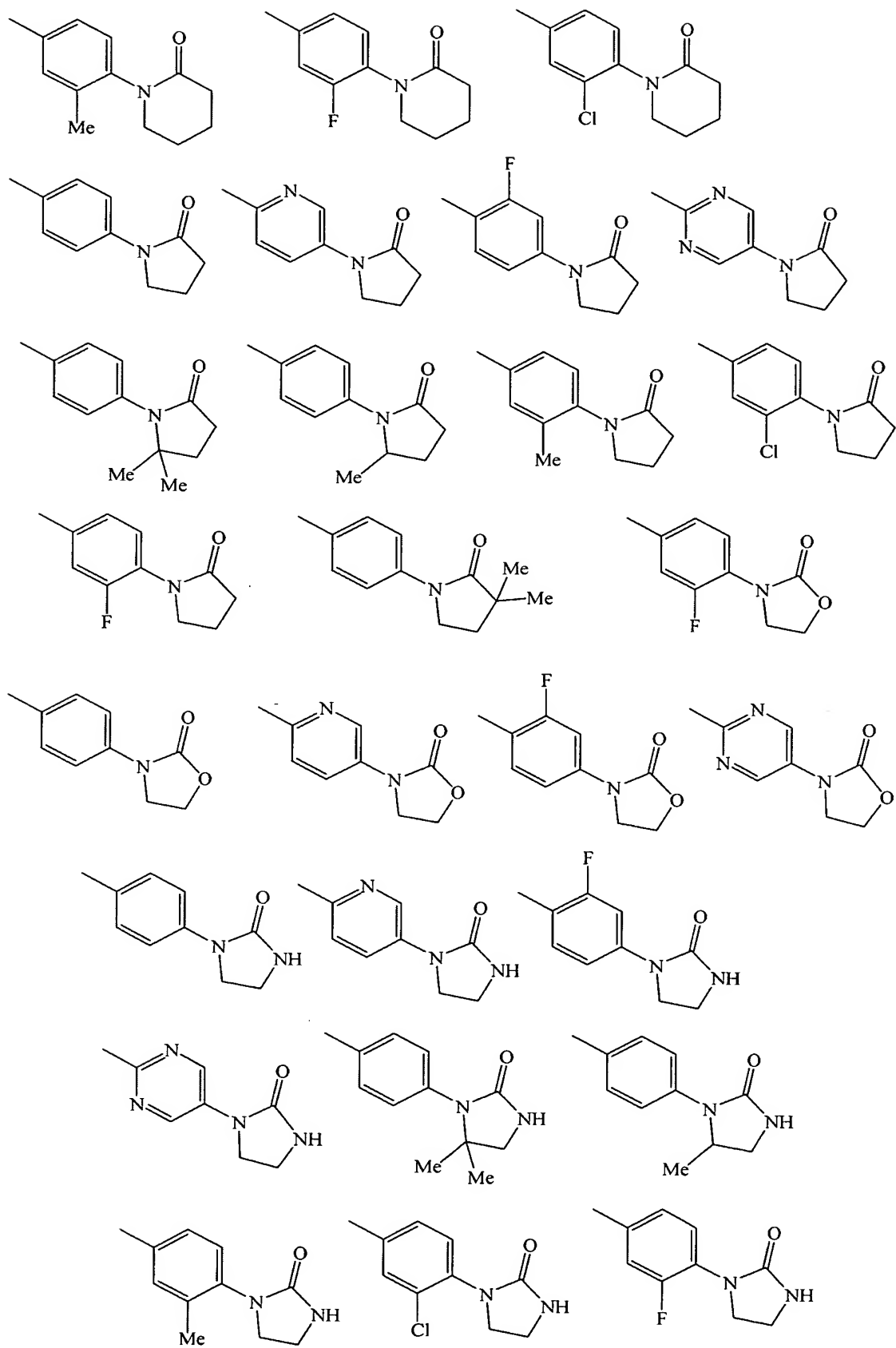
G is selected from:

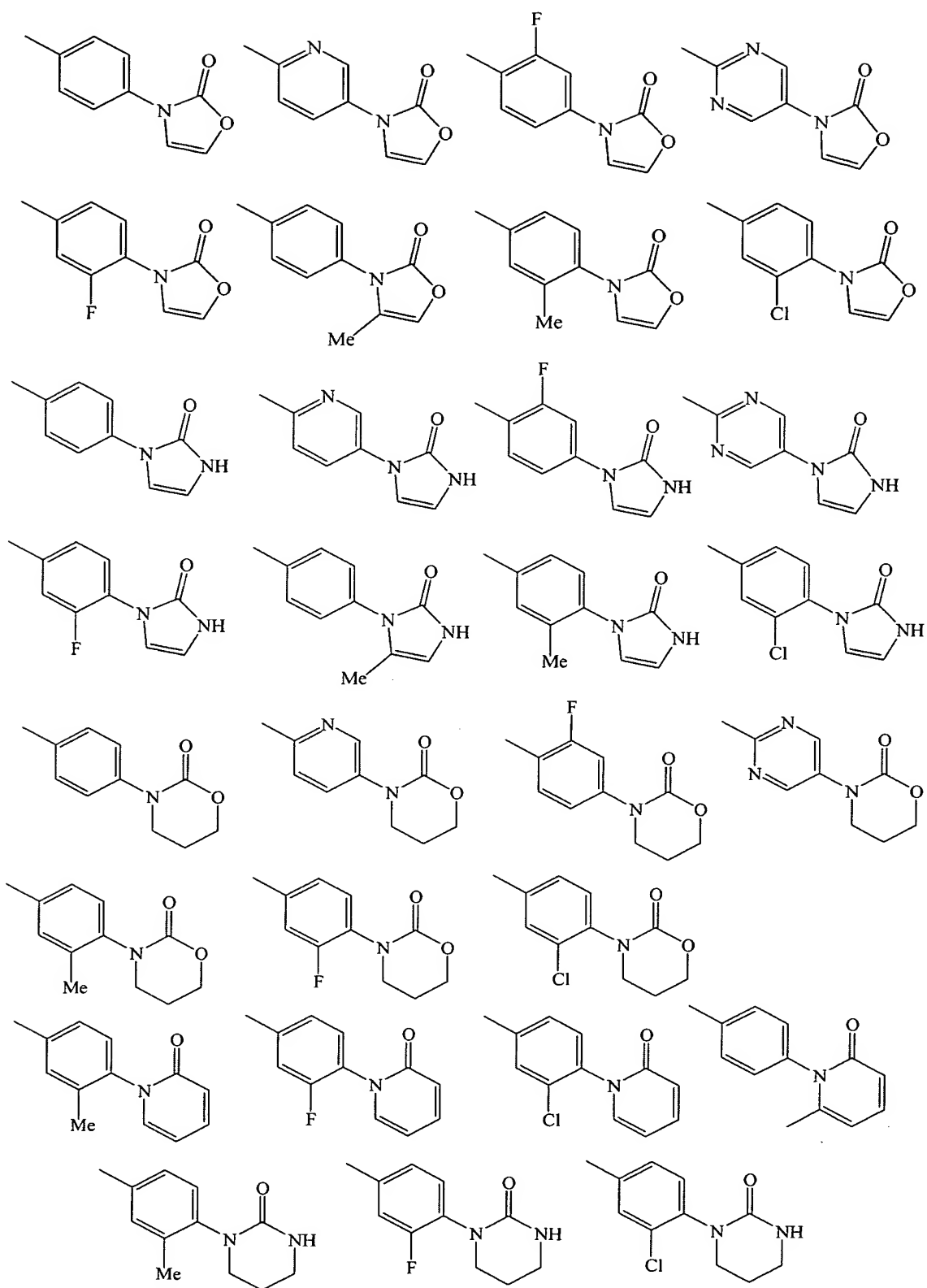


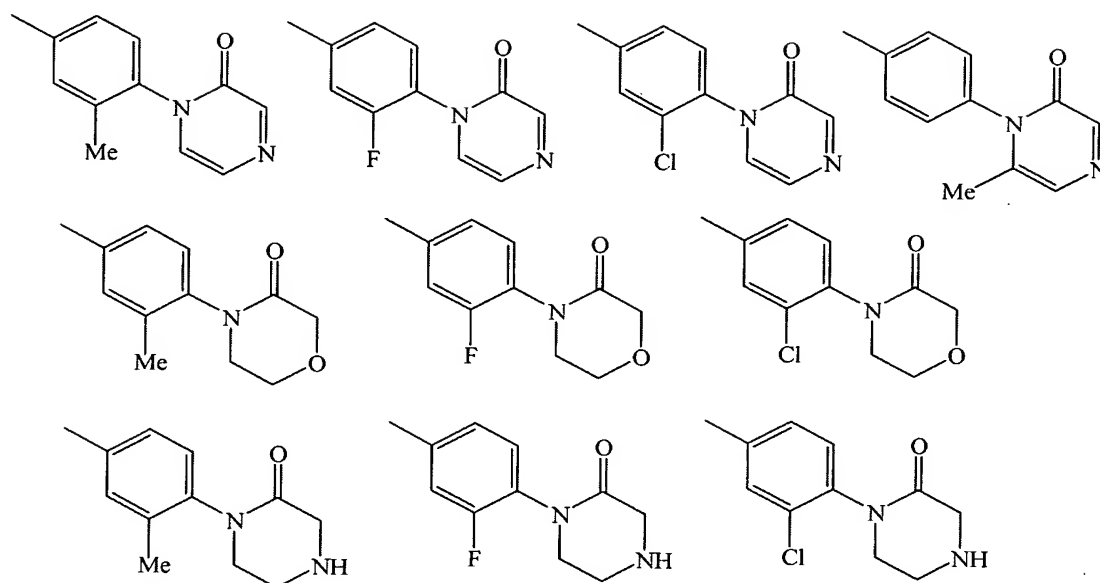
A-B is selected from:











G_1 is selected from $CH=CH$, $CH_2C(O)$, $C(O)CH_2$, NH , $C(O)NH$, $NHC(O)$, CH_2S , SCH_2 , $CH_2S(O)$, CH_2SO_2 , SO_2NH , $NHSO_2$, $NHCH_2C(O)NH$, $NHC(O)C(O)NH$, $NHC(O)C(S)NH$, and $NHC(S)C(O)NH$ and the right side of G_1 is attached to ring G, provided that Z does not form a N-S, NCH₂N, NCH₂O, or NCH₂S bond with either group to which it is attached;

Z is selected from $CH_2C(O)$, $C(O)CH_2$, NH , $C(O)NH$, $NHC(O)$, CH_2S , SCH_2 , $CH_2S(O)$, CH_2SO_2 , SO_2NH , and $NHSO_2$ and the right side of Z is attached to A, provided that Z does not form a N-S, NCH₂N, NCH₂O, or NCH₂S bond with either group to which it is attached;

R^{1a} is selected from H, R^{1b} , $C(CH_3)_2R^{1b}$, CH_2R^{1b} , and $CH_2CH_2R^{1b}$, provided that R^{1a} forms other than an N-halo, N-S, or N-CN bond;

R^{1b} is selected from CH_3 , CH_2CH_3 , F, Cl, Br, -CN, CF_3 , OR^2 , NR^2R^{2a} , $C(O)R^{2b}$, CO_2R^{2b} , CO_2R^{2a} , $S(O)_pR^{2b}$, $C(O)NR^2R^{2a}$, $C(O)NR^2R^{2b}$, $C(S)NR^2R^{2a}$, $SO_2NR^2R^{2a}$, $NR^2SO_2R^2$, cyclopropyl substituted with 0-2 R^{4b} , and 5-6 membered aromatic heterocycle consisting of carbon atoms and from 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{4b} , provided that R^{1b} forms other than an O-O, N-halo, N-S, or N-CN bond;

R^2 , at each occurrence, is selected from H, CH_3 , CH_2CH_3 , $CH_2CH_2CH_3$, $CH(CH_3)_2$, phenyl substituted with 0-1 R^{4b} , benzyl substituted with 0-1 R^{4b} , and 5-6

membered aromatic heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-1 R^{4b};

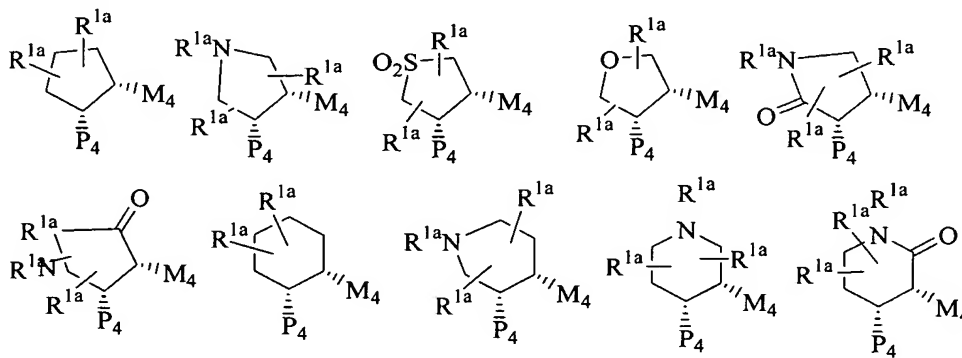
- R^{2a}, at each occurrence, is selected from H, CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, C(CH₃)₃, cyclopropyl, benzyl substituted with 0-1 R^{4b}, phenyl substituted with 0-1 R^{4b}, and 5-6 membered aromatic heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-1 R^{4b};

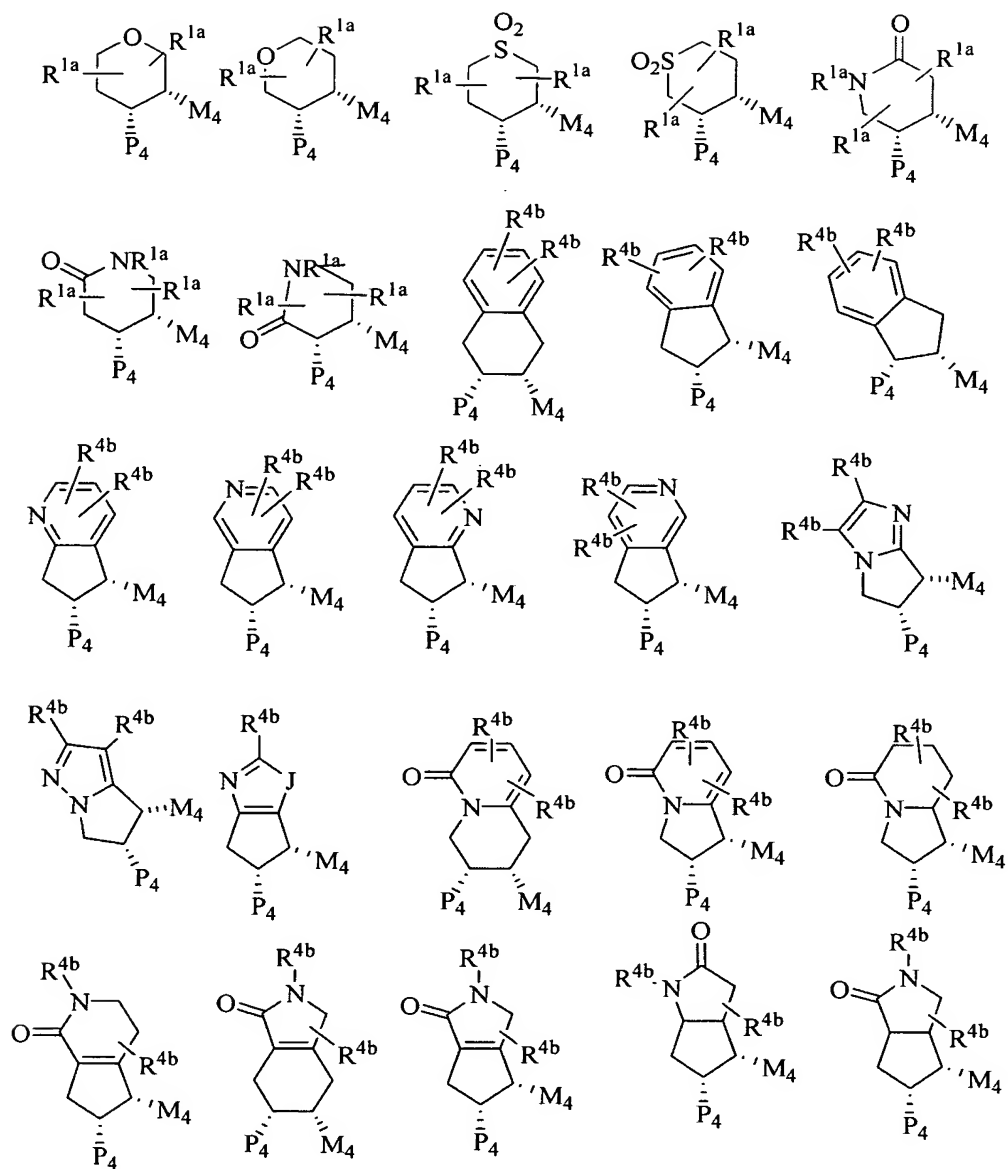
- alternatively, NR²R^{2a} forms a 4, 5, or 6 membered saturated, partially saturated, or unsaturated ring substituted with 0-1 R^{4b} and consisting of: carbon atoms, the nitrogen atom to which R² and R^{2a} are attached, and 0-1 additional heteroatoms selected from the group consisting of N, O, and S(O)_p;

- R^{2b}, at each occurrence, is selected from CF₃, OH, OCH₃, OCH₂CH₃, OCH₂CH₂CH₃, OCH(CH₃)₂, C₁₋₅ alkyl substituted with 0-3 R^{4b}, C₃₋₅ cycloalkyl substituted with 0-1 R^{4b}, benzyl substituted with 0-1 R^{4b}, phenyl substituted with 0-1 R^{4b}, and 5-6 membered heterocycle consisting of: carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and S(O)_p, and substituted with 0-1 R^{4b}; and

- R^{4b}, at each occurrence, is selected from H, =O, OR³, CH₂OR³, F, Cl, CH₃, CH₂CH₃, CH₂CH₂CH₃, CH(CH₃)₂, -CN, NO₂, NR³R^{3a}, CH₂NR³R^{3a}, C(O)R³, C(O)OR^{3c}, NR³C(O)R^{3a}, C(O)NR³R^{3a}, SO₂NR³R^{3a}, NR³SO₂-C₁₋₄ alkyl, NR³SO₂-phenyl, S(O)_p-C₁₋₄ alkyl, S(O)_p-phenyl, and CF₃.

5. A compound according to Claim 4, wherein the compound is selected from:



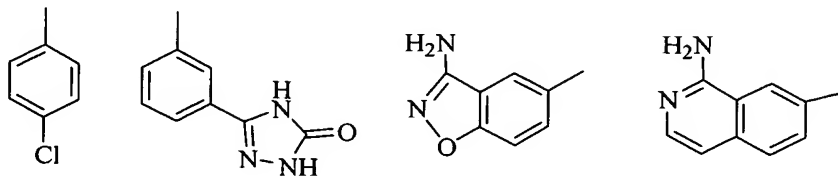


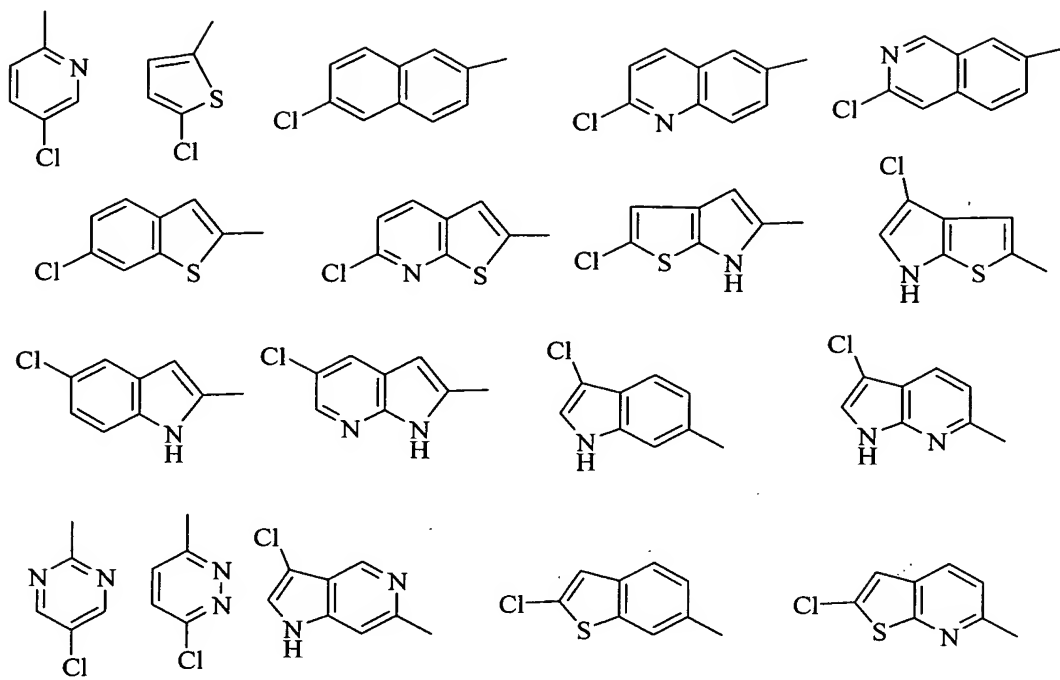
5

P_4 is G_1 -G;

M₄ is Z-A-B;

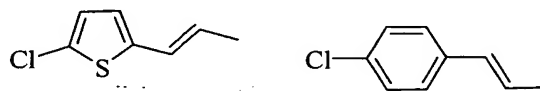
G is selected from:





G_1 is NHCO or NHC(=O)CONH;

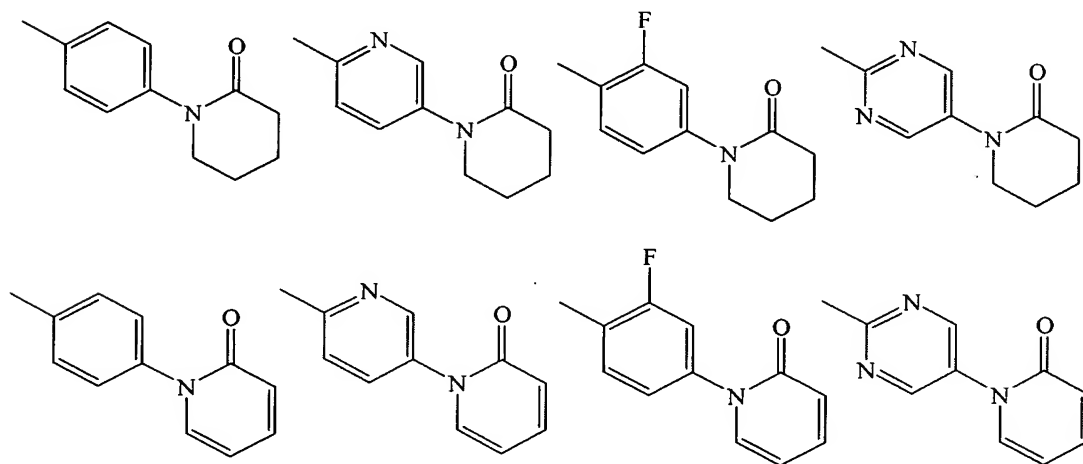
alternatively, G-G₁-is selected from:

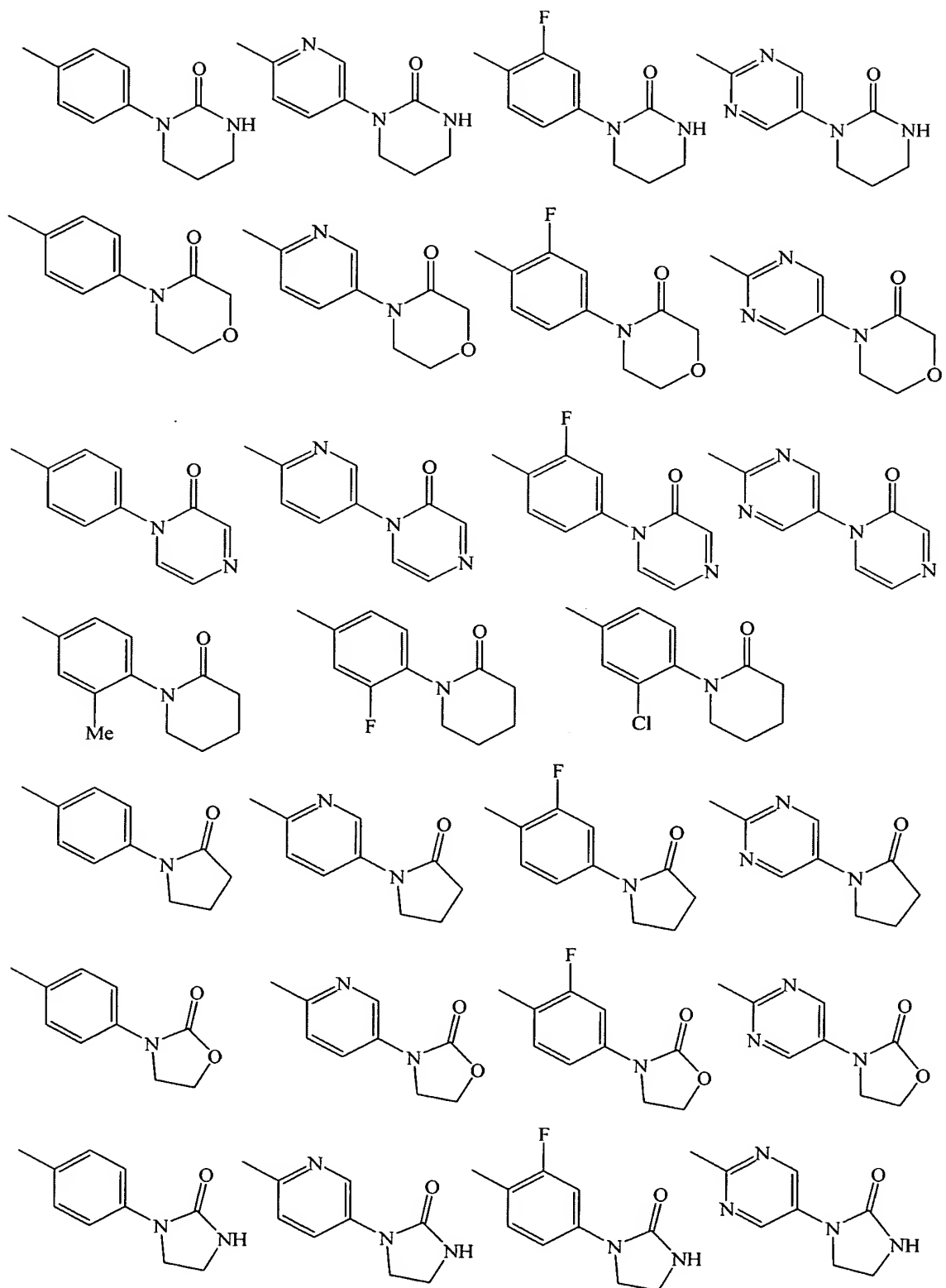


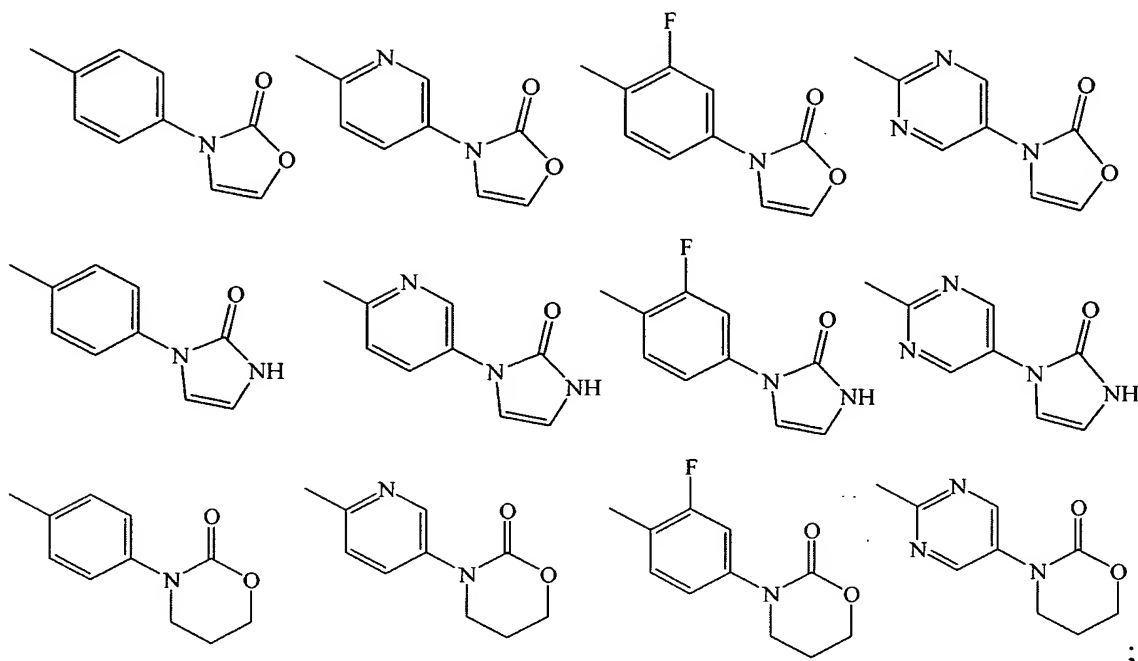
5

Z is NHCO or CONH;

A-B is selected from:



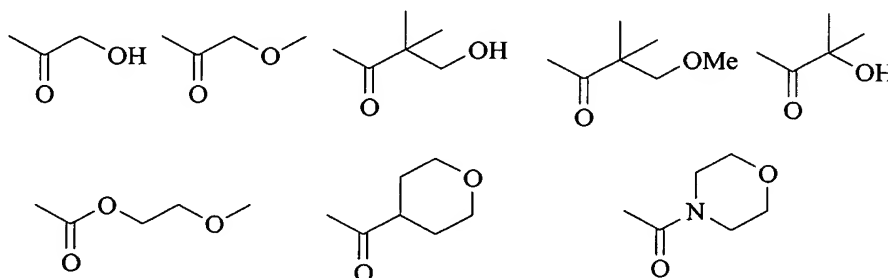


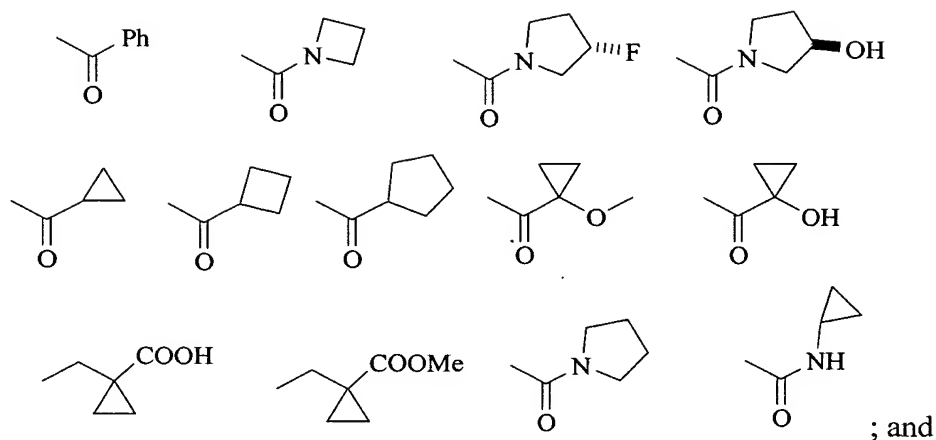


R^{1a} is selected from H, CH₃, CH₂CH₃, CH₂CH₂CH₃, OCH₃, CH₂OH,

- 5 CH₂CH₂OH, C(CH₃)₂OH, CH₂OCH₃, CH₂CH₂OCH₃, NH₂, CH₂NH₂, NHCH₃, CH₂NHCH₃, N(CH₃)₂, CH₂N(CH₃)₂, CO₂H, CH₂CO₂H, CH₂CH₂CO₂H, CH₂CH₂CO₂CH₂CH₃, COCH₃, COCH₂C(CH₃)₃, COCF₃, CO₂CH₃, CO₂CH₂CH₃, CO₂CH(CH₃)₂, CO₂C(CH₃)₃, CH₂CO₂CH₃, CH₂CH₂CO₂CH₂CH₃, S(O)₂CH₃, CH₂S(O)₂CH₃, C(O)NH₂, CONH(CH₃), CONH(CH₂CH₃), CONHC(CH₃)₃,
 10 CON(CH₃)₂, CON(CH₃)(CH₂CH₃), CON(CH₃)CH(CH₃)₂, CH₂C(O)NH₂, CH₂CON(CH₃)₂, CSN(CH₃)₂, SO₂NH₂, CH₂SO₂NH₂, NHSO₂CH₃, CH₂NHSO₂CH₃, SO₂CH₂CH₃, SO₂CH(CH₃)₂, SO₂CH₂CH₂CH₃, SO₂CH₂CH(CH₃)₂, SO₂CH₂CH₂OH, SO₂CH₂CH₂OCH₃, SO₂Ph, SO₂CH₂CF₃, and SO₂CF₂CF₃;

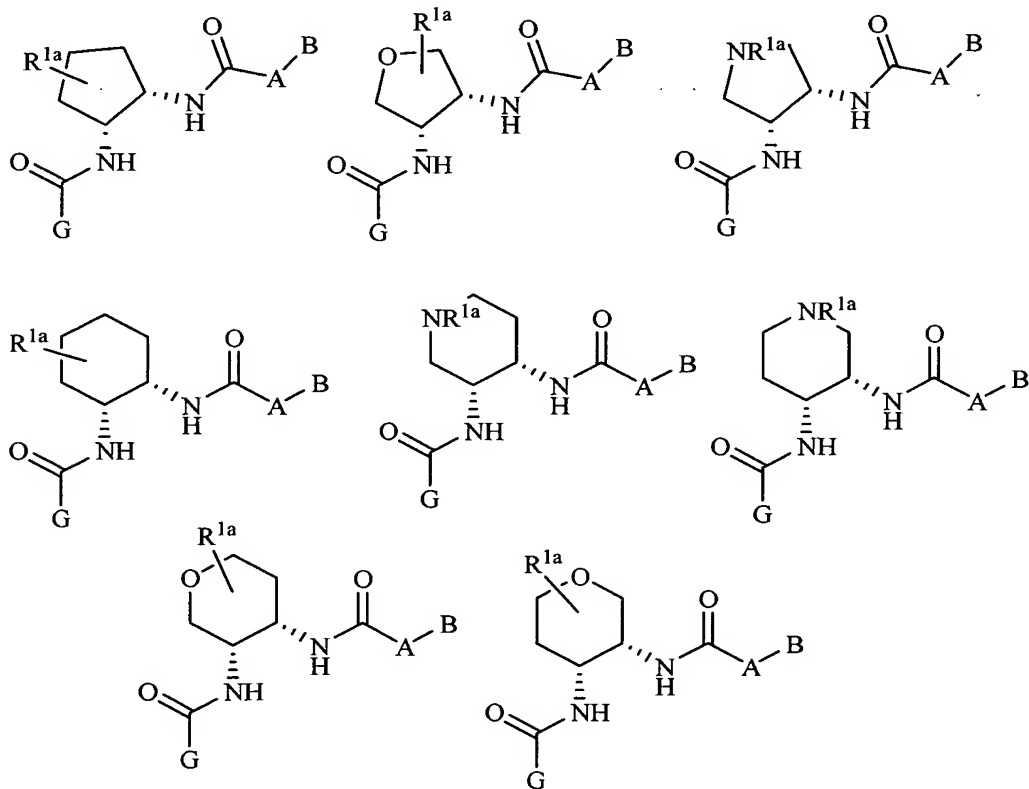
- 15 alternatively, R^{1a} is selected from:

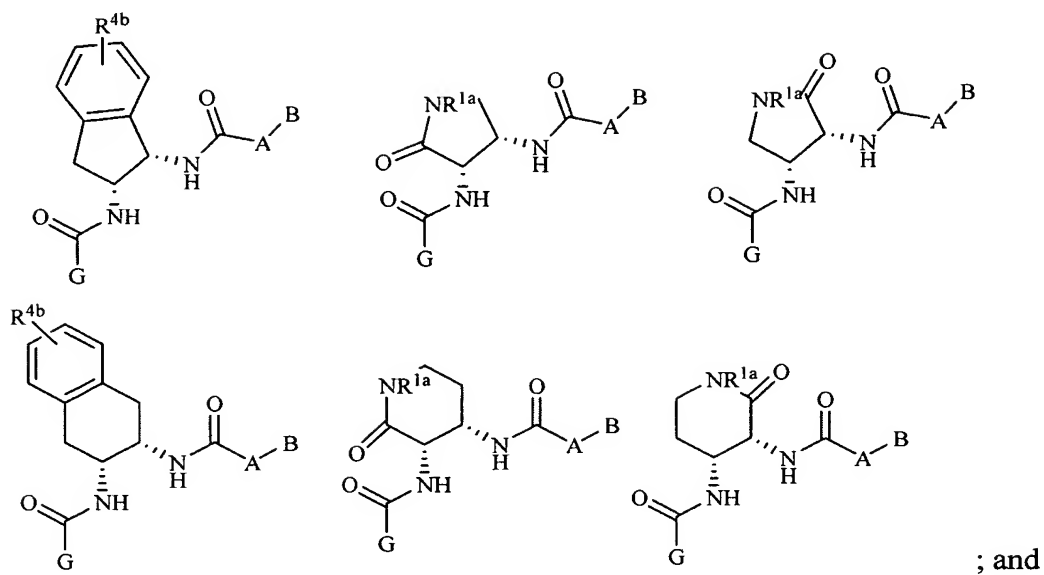




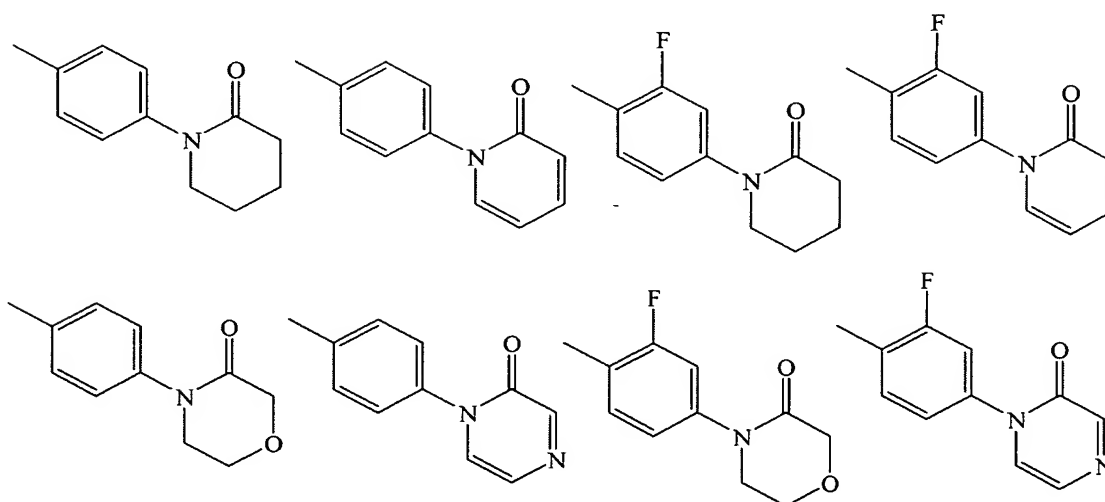
- R^{4b} , at each occurrence, is selected from H, =O, OR^3 , CH_2OR^3 , F, Cl, CH_3 , CH_2CH_3 , NR^3R^{3a} , $CH_2NR^3R^{3a}$, $C(O)R^3$, $C(O)OR^{3c}$, $NR^3C(O)R^{3a}$, $C(O)NR^3R^{3a}$,
 5 $SO_2NR^3R^{3a}$, NR^3SO_2 -phenyl, $S(O)_2CH_3$, $S(O)_2$ -phenyl, and CF_3 .

6. A compound according to Claim 5, wherein the compound is selected from:



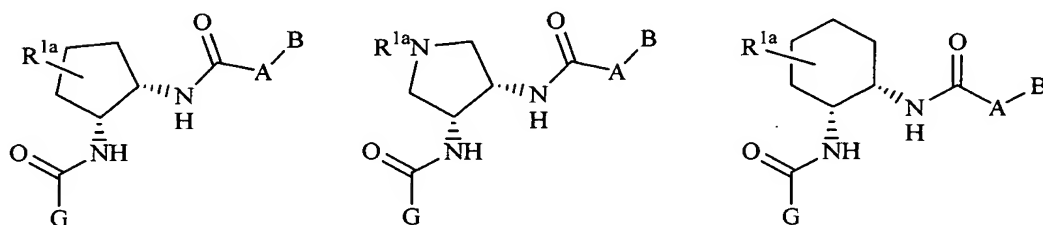


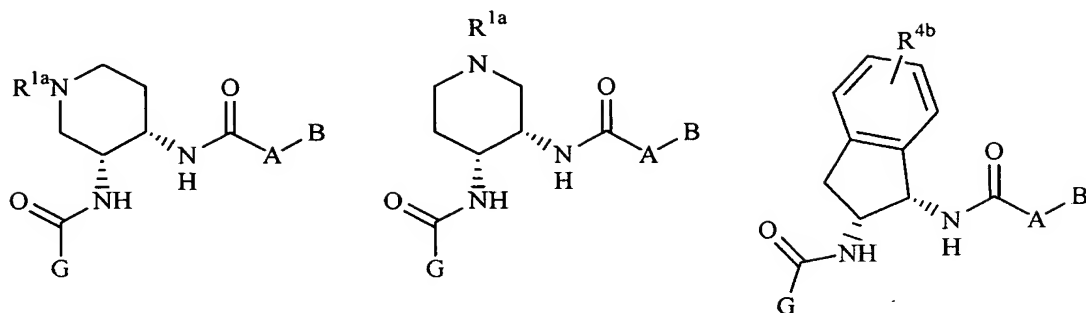
A-B is selected from:



5

7. A compound according to Claim 6, wherein the compound is selected from:





8. A compound according to Claim 1, wherein the compound is selected from the group:

- 5 (1*R*, 2*S*)-5-chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*S*, 2*R*)-5-chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-3-chloro-1H-indole-6-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-
- 10 benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-5-chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-piperidin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-3-chloro-1H-indole-6-carboxylic acid {2-[4-(2-oxo-piperidin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- 15 (1*S*, 2*R*)-4-methoxy-phenyl-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*S*, 2*R*)-5-chloro-1H-indole-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-5-chloro-1H-indole-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-
- 20 benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-5-chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclohexyl}-amide;
- (1*R*, 2*S*)-3-chloro-1H-indole-6-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclohexyl}-amide;
- 25 (1*R*, 2*S*)-3-chloro-1H-indole-6-carboxylic acid {2-[4-(2-oxo-piperidin-1-yl)-benzoylamino]-cyclohexyl}-amide;

- (1*R*, 2*S*)-*N*-(5-chloro-pyridin-2-yl)-*N'*-{2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-oxalamide;
- (1*S*, 2*S*)-3-chloro-1H-indole-6-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclohexyl}-amide;
- 5 (1*R*, 2*R*)-3-chloro-1H-indole-6-carboxylic acid {2-[4-(1-pyrrolidin-1-ylmethyl-cyclopropyl)-benzoylamino]-cyclohexyl}-amide;
- (1*S*, 3*R*, 4*S*)-3-[(3-chloro-1H-indole-6-carbonyl)-amino]-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentanecarboxylic acid methyl ester;
- (1*S*, 3*R*, 4*S*)-3-[(5-chloro-thiophene-2-carbonyl)-amino]-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentanecarboxylic acid methyl ester;
- 10 (1*R*, 2*S*, 4*S*)-5-chloro-thiophene-2-carboxylic acid {4-(2-methoxy-ethylcarbamoyl)-2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-3-chloro-1H-indole-6-carboxylic acid {2-[4-(3-oxo-morpholin-4-yl)-benzoylamino]-cyclohexyl}-amide;
- 15 (1*R*, 2*S*)-3-chloro-1H-indole-6-carboxylic acid {2-[4-(3-oxo-morpholin-4-yl)-benzoylamino]-cyclopentyl}-amide;
- cis*-3-chloro-1H-indole-6-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-phenylcarbamoyl]-cyclohexyl}-amide;
- (1*R*, 2*S*)-3-chloro-1H-indole-6-carboxylic acid {2-[3-methyl-4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclohexyl}-amide;
- 20 (1*R*, 2*S*)-3-chloro-1H-indole-6-carboxylic acid {2-[3-methyl-4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-3-chloro-1H-indole-6-carboxylic acid {2-[4-(2-oxo-azepan-1-yl)-benzoylamino]-cyclohexyl}-amide;
- 25 (1*R*, 2*S*)-3-chloro-1H-indole-6-carboxylic acid {2-[4-(2-oxo-azepan-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-*N*-[2-(6-chloro-naphthalene-2-sulfonylamino)-cyclohexyl]-4-(2-oxo-2H-pyridin-1-yl)-benzamide;
- (1*R*, 2*S*)-3-chloro-1H-indole-6-carboxylic acid {2-[4-(2-oxo-pyrrolidin-1-yl)-benzoylamino]-cyclohexyl}-amide;
- 30 (1*R*, 2*S*)-*N*-(5-chloro-pyridin-2-yl)-*N'*-{2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclohexyl}-oxalamide;

- (1*R*, 2*S*)-3-chloro-1*H*-indole-6-carboxylic acid {2-[4-(1,1-dioxo-1 λ ⁶-isothiazolidin-2-yl)-benzoylamino]-cyclohexyl}-amide;
- (1*R*, 2*S*)-3-chloro-1*H*-indole-6-carboxylic acid {2-[4-(2-oxo-tetrahydro-pyrimidin-1-yl)-benzoylamino]-cyclohexyl}-amide;
- 5 (1*R*, 2*S*)-3-chloro-1*H*-indole-6-carboxylic acid {2-[4-(2-oxo-[1,3]oxazinan-3-yl)-benzoylamino]-cyclohexyl}-amide;
- (1*R*, 2*S*)-3-chloro-1*H*-indole-6-carboxylic acid {2-[2-fluoro-4-(2-oxo-2*H*-pyridin-1-yl)-benzoylamino]-cyclohexyl}-amide;
- (1*R*, 2*S*)-4-chloro-phenylcarboxylic acid {2-[4-(2-oxo-2*H*-pyridin-1-yl)-
- 10 benzoylamino]-cyclohexyl}-amide;
- (1*R*, 2*S*)-3-chloro-1*H*-indole-6-carboxylic acid {2-[4-(2-oxo-2*H*-pyridin-1-yl)-benzoylamino]-indan-1-yl}-amide;
- (1*R*, 2*S*)-5-chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-2*H*-pyridin-1-yl)-benzoylamino]-indan-1-yl}-amide;
- 15 (1*S*, 2*R*)-5-chloro-thiophene-2-carboxylic acid {1-[4-(2-oxo-2*H*-pyridin-1-yl)-benzoylamino]-indan-2-yl}-amide;
- (1*S*, 2*R*)-3-chloro-1*H*-indole-6-carboxylic acid {1-[4-(2-oxo-2*H*-pyridin-1-yl)-benzoylamino]-indan-2-yl}-amide;
- (1*S*, 2*R*)-3-chloro-1*H*-indole-6-carboxylic acid {2-[4-(2-oxo-2*H*-pyridin-1-yl)-
- 20 benzoylamino]-indan-1-yl}-amide;
- (1*S*, 2*R*)-5-chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-2*H*-pyridin-1-yl)-benzoylamino]-indan-1-yl}-amide;
- (1*R*, 2*S*)-3-chloro-1*H*-indole-6-carboxylic acid {1-[4-(2-oxo-2*H*-pyridin-1-yl)-benzoylamino]-indan-2-yl}-amide;
- 25 *cis*-3-chloro-1*H*-indole-6-carboxylic acid {2-[4-(2-oxo-2*H*-pyridin-1-yl)-benzenesulfonylmethyl]-cyclohexyl}-amide;
- cis*-5-chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-2*H*-pyridin-1-yl)-benzenesulfonylmethyl]-cyclohexyl}-amide;
- cis*-*N*-[2-(4-chloro-benzenesulfonylmethyl)-cyclohexyl]-4-(2-oxo-2*H*-pyridin-
- 30 1-yl)-benzamide;
- cis*-*N*-[2-(4-chloro-benzenesulfonylmethyl)-cyclohexyl]-4-(2-oxo-2*H*-pyridin-1-yl)-benzamide;

- cis*-5-chloro-thiophene-2-sulfonic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzenesulfonylmethyl]-cyclohexyl}-amide;
- cis*-1-(4-chloro-phenyl)-3-{2-[4-(2-oxo-2H-pyridin-1-yl)-benzenesulfonylmethyl]-cyclohexyl}-urea;
- 5 *cis*-3-chloro-1H-indole-6-carboxylic acid {1-cyclopropanecarbonyl-3-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-4-yl}-amide;
- cis*-5-chloro-thiophene-2-carboxylic acid {1-cyclopropanecarbonyl-3-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-4-yl}-amide;
- 10 *cis*-1H-indole-6-carboxylic acid {1-cyclopropanecarbonyl-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-3-yl}-amide;
- (1*R*, 2*S*)-4-chloro-phenylcarboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-4-chloro-3-fluorophenylcarboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- 15 (1*R*, 2*S*)-4-chloro-3-methylphenylcarboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-4-chloro-3-methoxyphenylcarboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-5-methyl-thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- 20 (1*R*, 2*S*)-1H-indole-6-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-6-chloro-naphthalene-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclohexyl}-amide;
- 25 (1*R*, 2*S*)-6-chloro-1H-indole-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-5-chloro-1H-indole-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-5-chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- 30 (1*R*, 2*S*)-3-chloro-1H-indole-6-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclohexyl}-amide;

- (1*R*, 2*S*)-N-{2-[(5-chloro-thiophen-2-yl)methyl]-amino}-cyclopentyl}-4-(2-oxo-2H-pyridin-1-yl)-benzamide;
- (1*R*, 2*S*)-3-chloro-1H-indole-6-carboxylic acid {2-[4-(2-oxo-2H-pyrazin-1-yl)-benzoylamino]-cyclohexyl}-amide;
- 5 (1*R*, 2*S*)-5-chloro-thiophene-2-carboxylic acid {1-[4-(2-oxo-2H-pyrazin-1-yl)-benzoylamino]-indan-2-yl}-amide;
- cis*-3-chloro-1H-indole-6-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzenesulfonylamino]-cyclohexyl}-amide;
- (1*R*, 2*S*, 4*S*)-5-chloro-thiophene-2-carboxylic acid {4-dimethylcarbamoyl-2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- 10 (1*R*, 2*S*, 4*S*)-5-chloro-thiophene-2-carboxylic acid {4-cyclopropylcarbamoyl-2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*, 4*S*)-5-chloro-thiophene-2-carboxylic acid {4-(morpholine-4-carbonyl)-2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- 15 *cis*-3-chloro-1H-indole-6-carboxylic acid {1-methyl-3-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-4-yl}-amide;
- cis*-3-chloro-1H-indole-6-carboxylic acid {1-isopropyl-3-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-4-yl}-amide;
- (1*R*, 2*S*)-6-chloro-benzo[b]thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- 20 (1*R*, 2*S*)-6-chloro-benzo[b]thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclohexyl}-amide;
- (1*R*, 2*S*)-5-chloro-thiophene-2-carboxylic acid {2-[2-fluoro-4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- 25 (1*R*, 2*S*)-5-chloro-thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyrazin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1*R*, 2*S*)-6-chloro-benzo[b]thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyrazin-1-yl)-benzoylamino]-cyclohexyl}-amide;
- (1*S*, 2*R*)-2-oxo-2H-[1,2']bipyridinyl-5'-carboxylic acid {2-[(5-chloro-thiophene-2-carbonyl)-amino]-cyclopentyl}-amide;
- 30 (1*R*, 2*S*)-5-chloro-thiophene-2-carboxylic acid {2-[3-methyl-4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;

- (3R,4S)-3-[(5-chloro-thiophene-2-carbonyl)-amino]-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidine-1-carboxylic acid 9H-fluoren-9-yl methyl ester;
- (3R,4S)-5-chloro-thiophene-2-carboxylic acid {4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidin-3-yl}-amide;
- 5 (3R,4S)-3-[(3-chloro-1H-indole-6-carbonyl)-amino]-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidine-1-carboxylic acid 9H-fluoren-9-ylmethyl ester;
- (3R,4S)-3-chloro-1H-indole-6-carboxylic acid {4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidin-3-yl}-amide;
- (3R,4S)-5-chloro-thiophene-2-carboxylic acid {1-acetyl-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidin-3-yl}-amide;
- 10 (3R,4S)-3-chloro-1H-indole-5-carboxylic acid {1-acetyl-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidin-3-yl}-amide;
- (3R,4S)-5-chloro-thiophene-2-carboxylic acid {1-cyclopropanecarbonyl-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidin-3-yl}-amide;
- 15 (3R,4S)-5-chloro-thiophene-2-carboxylic acid {1-(2,2-dimethyl-propionyl)-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidin-3-yl}-amide;
- (3R,4S)-5-chloro-thiophene-2-carboxylic acid {4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-1-propionyl-pyrrolidin-3-yl}-amide;
- (3R,4S)-5-chloro-thiophene-2-carboxylic acid {1-(2-methoxy-acetyl)-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidin-3-yl}-amide;
- 20 (3R,4S)-5-chloro-thiophene-2-carboxylic acid {1-isobutyryl-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidin-3-yl}-amide;
- (3R,4S)-5-chloro-thiophene-2-carboxylic acid {1-benzoyl-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidin-3-yl}-amide;
- 25 (3R,4S)-5-chloro-thiophene-2-carboxylic acid {1-methanesulfonyl-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidin-3-yl}-amide;
- (3R,4S)-5-chloro-thiophene-2-carboxylic acid {1-ethanesulfonyl-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidin-3-yl}-amide;
- (3R,4S)-5-chloro-thiophene-2-carboxylic acid [4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-1-(propane-2-sulfonyl)-pyrrolidin-3-yl]-amide;
- 30 (3R,4S)-5-chloro-thiophene-2-carboxylic acid [4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-1-(pyrrolidine-1-carbonyl)-pyrrolidin-3-yl]-amide;

- (3R,4S)-3-[(3-chloro-1H-indole-5-carbonyl)-amino]-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidine-1-carboxylic acid ethyl ester;
- (3R,4S)-3-[(5-chloro-thiophene-2-carbonyl)-amino]-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidine-1-carboxylic acid methyl ester;
- 5 (3R,4S)-3-[(5-chloro-thiophene-2-carbonyl)-amino]-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidine-1-carboxylic acid ethyl ester;
- (3R,4S)-3-[(5-chloro-thiophene-2-carbonyl)-amino]-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-pyrrolidine-1-carboxylic acid 2-methoxy-ethyl ester;
- (1S,3R,4S)-3-[(5-chloro-thiophene-2-carbonyl)-amino]-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentanecarboxylic acid;
- 10 (1R,2S,4S)-5-chloro-thiophene-2-carboxylic acid {4-hydroxymethyl-2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1S,2R)-5-chloro-thiophene-2-carboxylic acid {1-[4-(2-oxo-piperidin-1-yl)-benzoylamino]-indan-2-yl}-amide;
- 15 (3S,4R)-5-chloro-thiophene-2-carboxylic acid {4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-tetrahydro-furan-3-yl}-amide;
- (3S,4R)-3-chloro-1H-indole-6-carboxylic acid {4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-tetrahydro-furan-3-yl}-amide;
- (1R,2S)-6-chloro-naphthalene-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- 20 (1R,2S)-5-chloro-3a,7a-dihydro-benzo[b]thiophene-2-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1R,2S)-3-chloro-phenyl-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- 25 (1R,2S)-2-chloro-phenyl-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1R,2S)-3,4-dichloro-phenyl-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1R,2S)-4-chloro-2-fluoro-phenyl-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- 30 (1R,2S)-2,4-dichloro-phenyl-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;

- (1R,2S)-4-chloro-2-methyl-phenyl-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1R,2S)-4-methoxy-phenyl-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- 5 (1R,2S)-3-methoxy-phenyl-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1R,2S)-2-chloro-thiazole-5-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- (1R,2S)-N-{2-[3-(4-chloro-phenyl)-ureido]-cyclopentyl}-4-(2-oxo-2H-pyridin-1-yl)-benzamide;
- 10 (1S,2R)-[2,2']bithiophenyl-5-carboxylic acid {2-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-cyclopentyl}-amide;
- cis*-3-[(3-chloro-1H-indole-6-carbonyl)-amino]-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidine-1-carboxylic acid *tert*-butyl ester;
- 15 *cis*-3-chloro-1H-indole-6-carboxylic acid {4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-3-yl}-amide;
- cis*-5-chloro-thiophene-2-carboxylic acid {1-cyclopropanecarbonyl-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-3-yl}-amide;
- cis*-3-chloro-1H-indole-6-carboxylic acid {1-methanesulfonyl-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-3-yl}-amide;
- 20 *cis*-3-chloro-1H-indole-6-carboxylic acid {1-acetyl-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-3-yl}-amide;
- cis*-3-[(3-chloro-1H-indole-6-carbonyl)-amino]-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidine-1-carboxylic acid ethyl ester;
- 25 *cis*-3-chloro-1H-indole-6-carboxylic acid {1-dimethylcarbamoyl-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-3-yl}-amide;
- cis*-3-chloro-1H-indole-6-carboxylic acid {1-cyclopropanecarbonyl-4-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-3-yl}-amide;
- cis*-4-[(3-chloro-1H-indole-6-carbonyl)-amino]-3-[4-(2-oxo-piperidin-1-yl)-benzoylamino]-piperidine-1-carboxylic acid *tert*-butyl ester;
- 30 *cis*-4-[(3-chloro-1H-indole-6-carbonyl)-amino]-3-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidine-1-carboxylic acid *tert*-butyl ester;

cis-3-chloro-1H-indole-6-carboxylic acid {1-methanesulfonyl-3-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-4-yl}-amide;

cis-3-chloro-1H-indole-6-carboxylic acid {1-acetyl-3-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-4-yl}-amide;

5 *cis*-3-chloro-1H-indole-6-carboxylic acid {1-dimethylcarbamoyl-3-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-4-yl}-amide;

cis-{1-cyclopropanecarbonyl-3-[4-(2-oxo-2H-pyridin-1-yl)-benzoylamino]-piperidin-4-yl}-carbamic acid benzyl ester;

(3R,4S)-3-(2-chlorothiophene-5-carboxamido)-N-methyl-4-(4-(2-oxopyridin-10 1(2H)-yl)benzamido)pyrrolidine-1-carboxamide; and

(3R,4S)-3-(2-chlorothiophene-5-carboxamido)-N,N-dimethyl-4-(4-(2-oxopyridin-1(2H)-yl)benzamido)pyrrolidine-1-carboxamide;
or a pharmaceutically acceptable salt form thereof.

15 9. A compound according to Claim 1, wherein the compound is selected from Examples 128-429 of Table 1.

10. A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound of Claim 1 or a
20 pharmaceutically acceptable salt form thereof.

11. A method for treating a thromboembolic disorder, comprising:
administering to a patient in need thereof a therapeutically effective amount of a
compound of Claim 1 or a pharmaceutically acceptable salt form thereof.

25 12. A method according to Claim 11, wherein the thromboembolic disorder is selected from the group consisting of arterial cardiovascular thromboembolic disorders, venous cardiovascular thromboembolic disorders, and thromboembolic disorders in the chambers of the heart.

30 13. A method according to Claim 11, wherein the thromboembolic disorder is selected from unstable angina, an acute coronary syndrome, first myocardial

infarction, recurrent myocardial infarction, ischemic sudden death, transient ischemic attack, stroke, atherosclerosis, peripheral occlusive arterial disease, venous thrombosis, deep vein thrombosis, thrombophlebitis, arterial embolism, coronary arterial thrombosis, cerebral arterial thrombosis, cerebral embolism, kidney embolism, pulmonary embolism, and thrombosis resulting from (a) prosthetic valves or other implants, (b) indwelling catheters, (c) stents, (d) cardiopulmonary bypass, (e) hemodialysis, or (f) other procedures in which blood is exposed to an artificial surface that promotes thrombosis.

- 10 14. A method for treating a thromboembolic disorder, comprising:
administering to a patient in need thereof a therapeutically effective amount of a first
and second therapeutic agent, wherein the first therapeutic agent is compound of
Claim 1 or a pharmaceutically acceptable salt thereof and the second therapeutic agent
is at least one agent selected from a second factor Xa inhibitor, an anti-coagulant
15 agent, an anti-platelet agent, a thrombin inhibiting agent, a thrombolytic agent, and a
fibrinolytic agent.